Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071725\
Data File : BF143147.D
Acq On : 17 Jul 2025 14:34
Operator : RC/JU
Sample : SSTDICC080
Misc :
ALS Vial : 10 Sample Multiplier: 1

Quant Time: Jul 17 15:05:57 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF071725.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 17 14:54:43 2025
Response via : Initial Calibration

Instrument :
BNA_F
ClientSampleId :
SSTDICC080

Manual Integrations APPROVED

Reviewed By :Rahul Chavli 07/18/2025
Supervised By :Jagrut Upadhyay 07/18/2025

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	6.963	152	145034	20.000	ng		0.00
21) Naphthalene-d8	8.245	136	544516	20.000	ng		0.00
39) Acenaphthene-d10	10.004	164	275619	20.000	ng		0.00
64) Phenanthrene-d10	11.486	188	429046	20.000	ng		0.00
76) Chrysene-d12	14.127	240	245351	20.000	ng		0.00
86) Perylene-d12	15.627	264	296687	20.000	ng		0.00
System Monitoring Compounds							
5) 2-Fluorophenol	5.587	112	1300344	141.880	ng		0.00
7) Phenol-d6	6.593	99	1639937	142.159	ng		0.00
23) Nitrobenzene-d5	7.534	82	1644581	150.551	ng		0.01
42) 2,4,6-Tribromophenol	10.792	330	437054	153.498	ng		0.00
45) 2-Fluorobiphenyl	9.322	172	2623071	130.157	ng		0.00
79) Terphenyl-d14	13.069	244	2202828	133.606	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	2.793	88	323993	74.710	ng		99
3) Pyridine	3.569	79	842446	74.844	ng		99
4) n-Nitrosodimethylamine	3.528	42	445612	76.677	ng		97
6) Aniline	6.628	93	1170943	72.831	ng		98
8) 2-Chlorophenol	6.751	128	699193	72.781	ng		99
9) Benzaldehyde	6.516	77	493544	57.055	ng		99
10) Phenol	6.610	94	888041m	70.663	ng		
11) bis(2-Chloroethyl)ether	6.698	93	697645	72.503	ng		100
12) 1,3-Dichlorobenzene	6.910	146	734699	70.400	ng		100
13) 1,4-Dichlorobenzene	6.987	146	739462	70.360	ng		99
14) 1,2-Dichlorobenzene	7.140	146	709300	70.958	ng		100
15) Benzyl Alcohol	7.104	79	661847	75.139	ng		99
16) 2,2'-oxybis(1-Chloropr...	7.240	45	1265589	70.902	ng		99
17) 2-Methylphenol	7.210	107	595477	73.719	ng		99
18) Hexachloroethane	7.481	117	266665	73.293	ng		99
19) n-Nitroso-di-n-propyla...	7.387	70	528820	71.452	ng		99
20) 3+4-Methylphenols	7.369	107	673906	68.763	ng	#	87
22) Acetophenone	7.375	105	890890	69.107	ng	#	95
24) Nitrobenzene	7.551	77	751964	73.836	ng		100
25) Isophorone	7.792	82	1453353	75.366	ng		99
26) 2-Nitrophenol	7.863	139	364951	82.568	ng		100
27) 2,4-Dimethylphenol	7.898	122	649144	72.050	ng		99
28) bis(2-Chloroethoxy)met...	7.992	93	834021	72.134	ng		99
29) 2,4-Dichlorophenol	8.104	162	558665	73.527	ng		99
30) 1,2,4-Trichlorobenzene	8.187	180	588333	71.673	ng		99
31) Naphthalene	8.269	128	1854387	69.150	ng		100
32) Benzoic acid	8.028	122	423625m	79.801	ng		
33) 4-Chloroaniline	8.310	127	787243	71.895	ng		98
34) Hexachlorobutadiene	8.392	225	368327	73.453	ng		99
35) Caprolactam	8.698	113	178271	77.644	ng		96
36) 4-Chloro-3-methylphenol	8.792	107	614048	74.354	ng		99
37) 2-Methylnaphthalene	8.957	142	1137176	69.687	ng		100
38) 1-Methylnaphthalene	9.057	142	1168702	69.550	ng		100
40) 1,2,4,5-Tetrachloroben...	9.128	216	568070	71.389	ng		99
41) Hexachlorocyclopentadiene	9.122	237	412756	79.719	ng		100
43) 2,4,6-Trichlorophenol	9.234	196	412148	74.838	ng		100

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Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
44) 2,4,5-Trichlorophenol	9.275	196	423565	76.888	ng		99
46) 1,1'-Biphenyl	9.428	154	1492094	69.388	ng		100
47) 2-Chloronaphthalene	9.451	162	1131626	71.122	ng		100
48) 2-Nitroaniline	9.534	65	404023	80.971	ng		96
49) Acenaphthylene	9.869	152	1868062	70.057	ng		100
50) Dimethylphthalate	9.728	163	1288559	71.723	ng		99
51) 2,6-Dinitrotoluene	9.781	165	292718	80.093	ng		99
52) Acenaphthene	10.039	154	1120315	71.086	ng		99
53) 3-Nitroaniline	9.951	138	333980	78.128	ng		98
54) 2,4-Dinitrophenol	10.051	184	133446	80.516	ng	#	40
55) Dibenzofuran	10.210	168	1615399	69.038	ng		99
56) 4-Nitrophenol	10.098	139	252698	79.161	ng		99
57) 2,4-Dinitrotoluene	10.186	165	376124	82.029	ng		96
58) Fluorene	10.551	166	1197731	68.203	ng		99
59) 2,3,4,6-Tetrachlorophenol	10.322	232	340118	74.533	ng		100
60) Diethylphthalate	10.422	149	1276217	71.870	ng		99
61) 4-Chlorophenyl-phenyle...	10.539	204	614962	70.326	ng		99
62) 4-Nitroaniline	10.569	138	291926	79.680	ng		98
63) Azobenzene	10.704	77	1318720	71.255	ng		99
65) 4,6-Dinitro-2-methylph...	10.592	198	186415	78.948	ng		98
66) n-Nitrosodiphenylamine	10.663	169	1087990	72.014	ng		98
67) 4-Bromophenyl-phenylether	11.033	248	382122	74.734	ng		98
68) Hexachlorobenzene	11.104	284	399107	74.680	ng		96
69) Atrazine	11.180	200	324126	77.819	ng		99
70) Pentachlorophenol	11.286	266	257176	81.829	ng		98
71) Phenanthrene	11.516	178	1602208	70.331	ng		100
72) Anthracene	11.563	178	1641264	70.640	ng		100
73) Carbazole	11.710	167	1435444	70.754	ng		99
74) Di-n-butylphthalate	12.045	149	1699503	74.034	ng		100
75) Fluoranthene	12.698	202	1498103	71.645	ng		99
78) Pyrene	12.927	202	1466205	70.021	ng		100
80) Butylbenzylphthalate	13.539	149	521631	81.353	ng		99
81) Benzo(a)anthracene	14.116	228	1212056	73.788	ng		100
82) 3,3'-Dichlorobenzidine	14.074	252	393448	71.867	ng		99
83) Chrysene	14.151	228	1105636	74.863	ng		100
84) Bis(2-ethylhexyl)phtha...	14.104	149	774366	79.790	ng		99
85) Di-n-octyl phthalate	14.716	149	1434055	81.519	ng		99
87) Indeno(1,2,3-cd)pyrene	17.174	276	1604969	76.718	ng		99
88) Benzo(b)fluoranthene	15.186	252	1438518	79.367	ng		100
89) Benzo(k)fluoranthene	15.216	252	1111726	68.737	ng		100
90) Benzo(a)pyrene	15.568	252	1269253	76.006	ng		99
91) Dibenzo(a,h)anthracene	17.198	278	1289789	75.746	ng		99
92) Benzo(g,h,i)perylene	17.645	276	1265778	76.846	ng		100

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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