Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111825\
Data File : BP026144.D
Acq On : 18 Nov 2025 15:12
Operator : RC/JU
Sample : SSTDICC005
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SSTDICC005

Quant Time: Nov 19 00:38:35 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 19 00:37:08 2025
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.672	152	261259	20.000	ng		0.00
21) Naphthalene-d8	10.437	136	1023584	20.000	ng		-0.01
39) Acenaphthene-d10	14.313	164	643620	20.000	ng		-0.01
64) Phenanthrene-d10	17.131	188	1346698	20.000	ng		0.00
76) Chrysene-d12	21.583	240	1560462	20.000	ng		0.01
86) Perylene-d12	24.918	264	1801943	20.000	ng		0.01
System Monitoring Compounds							
5) 2-Fluorophenol	5.290	112	155849	9.842	ng		0.00
7) Phenol-d6	6.849	99	191404	9.528	ng		0.00
23) Nitrobenzene-d5	8.807	82	171702	10.379	ng		0.00
42) 2,4,6-Tribromophenol	15.831	330	77590	11.005	ng		0.00
45) 2-Fluorobiphenyl	12.919	172	460038	10.313	ng		-0.01
79) Terphenyl-d14	19.872	244	798858	10.439	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.220	88	33093	4.981	ng		96
3) Pyridine	3.620	79	91499	4.844	ng		97
6) Aniline	7.002	93	127562	4.764	ng		98
8) 2-Chlorophenol	7.243	128	86099	4.908	ng		99
10) Phenol	6.872	94	102514	4.615	ng		96
11) bis(2-Chloroethyl)ether	7.096	93	81759	4.649	ng		99
12) 1,3-Dichlorobenzene	7.561	146	98674	4.925	ng		99
13) 1,4-Dichlorobenzene	7.708	146	97655	4.815	ng		99
14) 1,2-Dichlorobenzene	8.019	146	94897	4.899	ng		98
16) 2,2'-oxybis(1-Chloropr...	8.190	45	115566	4.387	ng		95
17) 2-Methylphenol	8.108	107	69679	4.771	ng		97
18) Hexachloroethane	8.743	117	34390	4.908	ng		97
19) n-Nitroso-di-n-propyla...	8.466	70	58743	4.718	ng		99
22) Acetophenone	8.478	105	127156	4.927	ng	#	99
24) Nitrobenzene	8.849	77	86731	5.006	ng		99
25) Isophorone	9.378	82	167606	4.787	ng		99
26) 2-Nitrophenol	9.555	139	38871	6.573	ng		98
27) 2,4-Dimethylphenol	9.625	122	78015	5.220	ng		98
28) bis(2-Chloroethoxy)met...	9.849	93	108845	4.925	ng		99
29) 2,4-Dichlorophenol	10.090	162	76304	4.893	ng		97
30) 1,2,4-Trichlorobenzene	10.302	180	84591	5.002	ng		94
31) Naphthalene	10.484	128	276491	4.976	ng		99
33) 4-Chloroaniline	10.590	127	109680	5.080	ng		100
34) Hexachlorobutadiene	10.784	225	54446	5.030	ng		99
36) 4-Chloro-3-methylphenol	11.731	107	80314	4.924	ng		96
37) 2-Methylnaphthalene	12.107	142	187351	4.822	ng		98
38) 1-Methylnaphthalene	12.337	142	187981	4.977	ng		99
40) 1,2,4,5-Tetrachloroben...	12.478	216	97644	4.903	ng		99
43) 2,4,6-Trichlorophenol	12.725	196	61137	5.131	ng		98
44) 2,4,5-Trichlorophenol	12.796	196	68937	5.093	ng		97
46) 1,1'-Biphenyl	13.131	154	241240	4.917	ng		97
47) 2-Chloronaphthalene	13.166	162	187850	4.870	ng		99
48) 2-Nitroaniline	13.378	65	46485	5.239	ng		90
49) Acenaphthylene	14.031	152	304102	5.334	ng		99
50) Dimethylphthalate	13.766	163	236591	5.077	ng		100
51) 2,6-Dinitrotoluene	13.872	165	37885	4.835	ng		91

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	Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
52)	Acenaphthene	14.378	154	183221	4.803	ng		98
53)	3-Nitroaniline	14.201	138	46080	4.861	ng	#	97
55)	Dibenzofuran	14.725	168	286273	4.928	ng		99
57)	2,4-Dinitrotoluene	14.678	165	58186	5.089	ng		97
58)	Fluorene	15.384	166	232855	5.102	ng		100
59)	2,3,4,6-Tetrachlorophenol	14.954	232	58158	5.425	ng		99
60)	Diethylphthalate	15.160	149	240172	5.123	ng		100
61)	4-Chlorophenyl-phenyle...	15.384	204	115474	5.076	ng		97
62)	4-Nitroaniline	15.395	138	50292	5.003	ng		97
63)	Azobenzene	15.678	77	193050	4.580	ng		99
66)	n-Nitrosodiphenylamine	15.595	169	201167	4.795	ng		99
67)	4-Bromophenyl-phenylether	16.301	248	69962	4.680	ng		98
68)	Hexachlorobenzene	16.413	284	81728	4.805	ng		97
69)	Atrazine	16.589	200	73147	5.086	ng		99
71)	Phenanthrene	17.178	178	379854	4.882	ng		100
72)	Anthracene	17.266	178	377623	4.865	ng		99
73)	Carbazole	17.554	167	364556	4.990	ng		100
74)	Di-n-butylphthalate	18.148	149	410422	4.968	ng		100
75)	Fluoranthene	19.266	202	465623	5.107	ng		100
78)	Pyrene	19.642	202	493981	4.740	ng		100
80)	Butylbenzylphthalate	20.613	149	212828	6.092	ng		100
81)	Benzo(a)anthracene	21.560	228	535793	5.028	ng		99
83)	Chrysene	21.630	228	506265	5.012	ng		99
84)	Bis(2-ethylhexyl)phtha...	21.524	149	328471	5.749	ng		99
87)	Indeno(1,2,3-cd)pyrene	28.677	276	649473	4.913	ng		99
88)	Benzo(b)fluoranthene	23.854	252	535938	4.848	ng		100
89)	Benzo(k)fluoranthene	23.930	252	543312	4.807	ng		99
90)	Benzo(a)pyrene	24.754	252	504836	4.939	ng		99
91)	Dibenzo(a,h)anthracene	28.777	278	530340	4.920	ng		99
92)	Benzo(g,h,i)perylene	29.853	276	529220	5.083	ng		99

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

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