Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120125\
Data File : BP026201.D
Acq On : 01 Dec 2025 17:51
Operator : RC/JU
Sample : Q3719-10MS
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-1125-15BMS

Quant Time: Dec 01 18:17:49 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 01:25:31 2025
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.655	152	233034	20.000	ng	-0.02
21) Naphthalene-d8	10.413	136	911777	20.000	ng	-0.04
39) Acenaphthene-d10	14.290	164	569978	20.000	ng	-0.04
64) Phenanthrene-d10	17.095	188	1129941	20.000	ng	-0.04
76) Chrysene-d12	21.536	240	1151085	20.000	ng	-0.04
86) Perylene-d12	24.830	264	1323407	20.000	ng	-0.08
System Monitoring Compounds						
5) 2-Fluorophenol	5.278	112	923920	63.638	ng	-0.01
7) Phenol-d6	6.837	99	1166768	63.126	ng	-0.02
23) Nitrobenzene-d5	8.790	82	684932	42.670	ng	-0.02
42) 2,4,6-Tribromophenol	15.807	330	481439	65.112	ng	-0.03
45) 2-Fluorobiphenyl	12.896	172	1626385	41.820	ng	-0.04
79) Terphenyl-d14	19.836	244	2507666	44.421	ng	-0.04
Target Compounds						Qvalue
2) 1,4-Dioxane	3.220	88	250774	42.585	ng	98
3) Pyridine	3.608	79	662225	38.913	ng	100
4) n-Nitrosodimethylamine	3.514	42	304533	47.880	ng	89
6) Aniline	6.990	93	630018	25.816	ng	99
8) 2-Chlorophenol	7.225	128	788736	47.727	ng	99
9) Benzaldehyde	6.802	77	314375	32.457	ng	100
10) Phenol	6.861	94	930117	46.713	ng	97
11) bis(2-Chloroethyl)ether	7.084	93	701792	45.015	ng	98
12) 1,3-Dichlorobenzene	7.543	146	855675	48.060	ng	99
13) 1,4-Dichlorobenzene	7.690	146	866519	48.300	ng	100
14) 1,2-Dichlorobenzene	8.002	146	834465	48.438	ng	100
15) Benzyl Alcohol	7.890	79	624651	46.131	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.172	45	998081	45.585	ng	100
17) 2-Methylphenol	8.096	107	629245	46.424	ng	98
18) Hexachloroethane	8.719	117	309058	47.346	ng	98
19) n-Nitroso-di-n-propyla...	8.449	70	505235	44.000	ng	97
20) 3+4-Methylphenols	8.413	107	853730	45.454	ng	100
22) Acetophenone	8.461	105	1096937	47.332	ng	99
24) Nitrobenzene	8.831	77	764580	47.089	ng	99
25) Isophorone	9.361	82	1458953	45.847	ng	99
26) 2-Nitrophenol	9.537	139	401702	48.892	ng	98
27) 2,4-Dimethylphenol	9.602	122	703951	47.516	ng	100
28) bis(2-Chloroethoxy)met...	9.831	93	922117	46.269	ng	100
29) 2,4-Dichlorophenol	10.066	162	733579	49.541	ng	99
30) 1,2,4-Trichlorobenzene	10.284	180	776015	51.465	ng	99
31) Naphthalene	10.466	128	2311162	46.902	ng	99
32) Benzoic acid	9.737	122	551835	44.231	ng	98
33) 4-Chloroaniline	10.572	127	426200	20.338	ng	100
34) Hexachlorobutadiene	10.760	225	484470	49.402	ng	99
35) Caprolactam	11.355	113	240085	45.280	ng	94
36) 4-Chloro-3-methylphenol	11.702	107	762537	48.452	ng	100
37) 2-Methylnaphthalene	12.078	142	1498324	42.890	ng	100
38) 1-Methylnaphthalene	12.307	142	1552007	45.453	ng	99
40) 1,2,4,5-Tetrachloroben...	12.454	216	872851	50.645	ng	99
41) Hexachlorocyclopentadiene	12.437	237	1071835	95.026	ng	100
43) 2,4,6-Trichlorophenol	12.690	196	594644	50.327	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.766	196	656295	49.761	ng	100
46) 1,1'-Biphenyl	13.101	154	2111302	49.193	ng	99
47) 2-Chloronaphthalene	13.137	162	1637408	48.516	ng	99
48) 2-Nitroaniline	13.343	65	461630	49.187	ng	95
49) Acenaphthylene	14.001	152	2739172	49.204	ng	100
50) Dimethylphthalate	13.743	163	2043526	48.002	ng	100
51) 2,6-Dinitrotoluene	13.849	165	457796	54.271	ng	99
52) Acenaphthene	14.354	154	1541992	47.265	ng	99
53) 3-Nitroaniline	14.184	138	302286	30.436	ng	99
54) 2,4-Dinitrophenol	14.396	184	543215	106.681	ng	97
55) Dibenzofuran	14.701	168	2439496	48.278	ng	99
56) 4-Nitrophenol	14.513	139	830971	92.485	ng	95
57) 2,4-Dinitrotoluene	14.660	165	637580	50.621	ng	98
58) Fluorene	15.360	166	1962637	49.141	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.931	232	586135	52.385	ng	99
60) Diethylphthalate	15.137	149	2241447	52.266	ng	100
61) 4-Chlorophenyl-phenyle...	15.360	204	976947	49.162	ng	99
62) 4-Nitroaniline	15.378	138	436714	42.344	ng	96
63) Azobenzene	15.660	77	1905150	52.687	ng	99
65) 4,6-Dinitro-2-methylph...	15.431	198	362269	50.467	ng	98
66) n-Nitrosodiphenylamine	15.578	169	1775615	50.799	ng	100
67) 4-Bromophenyl-phenylether	16.278	248	634525	50.706	ng	100
68) Hexachlorobenzene	16.390	284	745760	51.358	ng	100
69) Atrazine	16.560	200	670034	50.835	ng	98
70) Pentachlorophenol	16.742	266	1025730	100.983	ng	100
71) Phenanthrene	17.137	178	3215634	50.519	ng	100
72) Anthracene	17.237	178	3257412	50.174	ng	100
73) Carbazole	17.513	167	2967469	48.125	ng	100
74) Di-n-butylphthalate	18.119	149	3754359	51.358	ng	100
75) Fluoranthene	19.231	202	3879717	50.166	ng	99
77) Benzidine	19.436	184	1205912	33.011	ng	99
78) Pyrene	19.607	202	3990288	52.870	ng	99
80) Butylbenzylphthalate	20.578	149	1665637	50.112	ng	99
81) Benzo(a)anthracene	21.519	228	3944904	49.901	ng	100
82) 3,3'-Dichlorobenzidine	21.442	252	952449	33.130	ng	99
83) Chrysene	21.583	228	3694023	49.581	ng	100
84) Bis(2-ethylhexyl)phtha...	21.477	149	2469683	49.085	ng	99
85) Di-n-octyl phthalate	22.742	149	4234005	48.125	ng	100
87) Indeno(1,2,3-cd)pyrene	28.577	276	5124135	51.628	ng	94
88) Benzo(b)fluoranthene	23.789	252	4141922	51.393	ng	100
89) Benzo(k)fluoranthene	23.854	252	4026385	50.019	ng	100
90) Benzo(a)pyrene	24.677	252	4013559	52.575	ng	99
91) Dibenzo(a,h)anthracene	28.642	278	4124078	50.760	ng	98
92) Benzo(g,h,i)perylene	29.718	276	4117869	50.992	ng	100

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

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