

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032421\
 Data File : BP005034.D
 Acq On : 24 Mar 2021 12:22
 Operator : CG/JU
 Sample : M1557-03DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PT-BN-WPDL

Quant Time: Mar 24 12:56:39 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 24 10:16:52 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	25867	20.00	ng	# 0.00
21) Naphthalene-d8	10.516	136	103948	20.00	ng	# 0.00
39) Acenaphthene-d10	14.381	164	69905	20.00	ng	0.00
64) Phenanthrene-d10	17.139	188	144442	20.00	ng	# 0.00
76) Chrysene-d12	21.233	240	147795	20.00	ng	0.00
86) Perylene-d12	23.521	264	149145	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.317	112	47017	27.97	ng	0.00
7) Phenol-d6	6.899	99	62575	25.99	ng	0.00
23) Nitrobenzene-d5	8.887	82	42863	17.69	ng	0.00
42) 2,4,6-Tribromophenol	15.881	330	19170	21.05	ng	0.01
45) 2-Fluorobiphenyl	12.998	172	93479	18.09	ng	0.00
79) Terphenyl-d14	19.710	244	153128	18.86	ng	0.00
Target Compounds						
						Qvalue
4) n-Nitrosodimethylamine	3.570	42	16688	25.64	ng	# 71
11) bis(2-Chloroethyl)ether	7.158	93	50575	28.00	ng	92
12) 1,3-Dichlorobenzene	7.616	146	19075	9.68	ng	# 81
13) 1,4-Dichlorobenzene	7.763	146	64014	32.01	ng	# 90
14) 1,2-Dichlorobenzene	8.075	146	32241	16.81	ng	# 89
18) Hexachloroethane	8.799	117	16911	22.19	ng	89
24) Nitrobenzene	8.928	77	20332	7.96	ng	# 78
30) 1,2,4-Trichlorobenzene	10.387	180	50278	25.13	ng	98
34) Hexachlorobutadiene	10.852	225	26924	20.91	ng	96
37) 2-Methylnaphthalene	12.187	142	73624	17.36	ng	# 94
41) Hexachlorocyclopentadiene	12.534	237	35683	28.82	ng	97
48) 2-Nitroaniline	13.457	65	27688	20.02	ng	# 65
50) Dimethylphthalate	13.840	163	82601	15.34	ng	99
51) 2,6-Dinitrotoluene	13.963	165	25492	20.09	ng	# 55
52) Acenaphthene	14.445	154	69482	16.73	ng	100
55) Dibenzofuran	14.781	168	112940	16.59	ng	98
57) 2,4-Dinitrotoluene	14.751	165	7530	4.49	ng	# 50
60) Diethylphthalate	15.210	149	110017	20.89	ng	100
61) 4-Chlorophenyl-phenyle...	15.434	204	49357	16.92	ng	95
62) 4-Nitroaniline	15.469	138	4093	3.41	ng	# 57
66) n-Nitrosodiphenylamine	15.645	169	63624	13.70	ng	96
67) 4-Bromophenyl-phenylether	16.328	248	39377	23.53	ng	94
71) Phenanthrene	17.181	178	217429	25.97	ng	99
74) Di-n-butylphthalate	18.110	149	155147	18.34	ng	# 97
78) Pyrene	19.510	202	226766	22.33	ng	98
83) Chrysene	21.269	228	110534	11.22	ng	98
84) Bis(2-ethylhexyl)phtha...	21.145	149	145444	26.32	ng	99
85) Di-n-octyl phthalate	22.033	149	74990	8.05	ng	# 93
87) Indeno(1,2,3-cd)pyrene	25.863	276	134863	12.06	ng	97
88) Benzo(b)fluoranthene	22.827	252	150058	14.17	ng	99
89) Benzo(k)fluoranthene	22.869	252	113073	11.14	ng	99
90) Benzo(a)pyrene	23.416	252	117351	12.36	ng	98
91) Dibenzo(a,h)anthracene	25.886	278	142584	14.74	ng	# 96
92) Benzo(g,h,i)perylene	26.592	276	171086	18.33	ng	# 93

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

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