

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080624\
 Data File : BN033273.D
 Acq On : 06 Aug 2024 14:55
 Operator : MA/JU
 Sample : PB162490BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB162490BSD

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 08/07/2024
 Supervised By :mohammad ahmed 08/08/2024

Quant Time: Aug 06 15:33:08 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN080524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 05 23:57:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.561	152	699m	0.400	ng	0.05
7) Naphthalene-d8	10.340	136	2758	0.400	ng	0.02
13) Acenaphthene-d10	14.137	164	1926	0.400	ng	0.01
19) Phenanthrene-d10	16.933	188	4480	0.400	ng	0.01
29) Chrysene-d12	21.134	240	4687	0.400	ng	# 0.00
35) Perylene-d12	23.309	264	5795	0.400	ng	0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.416	112	742m	0.429	ng	0.12
5) Phenol-d6	7.048	99	1003m	0.422	ng	0.06
8) Nitrobenzene-d5	8.984	82	772	0.348	ng	0.06
11) 2-Methylnaphthalene-d10	11.950	152	1501	0.367	ng	0.02
14) 2,4,6-Tribromophenol	15.729	330	299	0.417	ng	0.01
15) 2-Fluorobiphenyl	12.822	172	2596	0.394	ng	0.02
27) Fluoranthene-d10	18.964	212	4942	0.394	ng	0.01
31) Terphenyl-d14	19.568	244	3289	0.406	ng	0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	3.047	88	482	0.387	ng	# 85
3) n-Nitrosodimethylamine	3.494	42	251m	0.294	ng	
6) bis(2-Chloroethyl)ether	7.134	93	1613m	0.805	ng	
9) Naphthalene	10.394	128	3019	0.403	ng	95
10) Hexachlorobutadiene	10.522	225	603	0.433	ng	# 98
12) 2-Methylnaphthalene	12.030	142	1496	0.335	ng	97
16) Acenaphthylene	13.880	152	2958	0.385	ng	99
17) Acenaphthene	14.201	154	2309	0.416	ng	99
18) Fluorene	15.227	166	3080	0.410	ng	99
20) 4,6-Dinitro-2-methylph...	15.604	198	90	0.354	ng	85
21) 4-Bromophenyl-phenylether	16.126	248	938	0.395	ng	97
22) Hexachlorobenzene	16.200	284	1034	0.417	ng	97
23) Atrazine	16.424	200	814	0.364	ng	95
24) Pentachlorophenol	16.647	266	280	0.337	ng	93
25) Phenanthrene	16.970	178	4622	0.384	ng	100
26) Anthracene	17.094	178	3452	0.342	ng	# 88
28) Fluoranthene	18.997	202	6285	0.380	ng	99
30) Pyrene	19.364	202	6699	0.409	ng	99
32) Benzo(a)anthracene	21.134	228	4405	0.326	ng	97
33) Chrysene	21.178	228	8223	0.453	ng	98
34) Bis(2-ethylhexyl)phtha...	21.008	149	1639	0.505	ng	# 89
36) Indeno(1,2,3-cd)pyrene	25.504	276	9137	0.349	ng	98
37) Benzo(b)fluoranthene	22.660	252	6426	0.366	ng	98
38) Benzo(k)fluoranthene	22.707	252	8817	0.401	ng	98
39) Benzo(a)pyrene	23.233	252	6890	0.390	ng	97
40) Dibenzo(a,h)anthracene	25.504	278	7145	0.345	ng	94
41) Benzo(g,h,i)perylene	26.174	276	8498	0.358	ng	98

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

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