

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080222\
 Data File : BN020940.D
 Acq On : 02 Aug 2022 13:16
 Operator : CG/JU
 Sample : SP5565
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SP5565

Quant Time: Aug 02 22:33:31 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 04:19:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.809	152	2534	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	8125	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.456	164	4987	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.209	188	11671	0.400	ng/ul	0.00
17) Chrysene-d12	21.424	240	10363	0.400	ng/ul	0.00
23) Perylene-d12	23.774	264	8263	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	2942	0.952	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.207	152	4678	0.390	ng/ul	0.00
18) Fluoranthene-d10	19.254	212	12783	0.403	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.244	88	14737	4.429	ng/ul#	82
5) Naphthalene	10.651	128	16809	0.757	ng/ul	98
7) 2-Methylnaphthalene	12.279	142	9559	0.707	ng/ul	99
8) 1-Methylnaphthalene	12.499	142	11054	0.740	ng/ul	100
10) Acenaphthylene	14.179	152	14503	0.711	ng/ul	98
11) Acenaphthene	14.521	153	12597	0.739	ng/ul	95
12) Fluorene	15.511	166	16171	0.824	ng/ul	98
14) Pentachlorophenol	16.876	266	2497	1.340	ng/ul	100
15) Phenanthrene	17.252	178	26633	0.673	ng/ul	99
16) Anthracene	17.349	178	21599	0.635	ng/ul	98
19) Fluoranthene	19.282	202	29504	0.633	ng/ul	99
20) Pyrene	19.644	202	30304	0.621	ng/ul	96
21) Benzo(a)anthracene	21.409	228	23418	0.622	ng/ul	100
22) Chrysene	21.462	228	28819	0.619	ng/ul	99
24) Benzo(b)fluoranthene	23.058	252	26964	0.684	ng/ul	97
25) Benzo(k)fluoranthene	23.104	252	28482	0.666	ng/ul	97
26) Benzo(a)pyrene	23.669	252	24118	0.648	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.192	276	28220	0.645	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.212	278	22779	0.636	ng/ul	95
29) Benzo(g,h,i)perylene	26.933	276	25670	0.645	ng/ul	98

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

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