

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080222\
 Data File : BN020965.D
 Acq On : 03 Aug 2022 05:25
 Operator : CG/JU
 Sample : PB146650BS
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS650

Quant Time: Aug 03 06:03:35 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	7.810	152	2206	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	7362	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.457	164	4093	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.210	188	8013	0.400	ng/ul	0.00
17) Chrysene-d12	21.415	240	6884	0.400	ng/ul	0.00
23) Perylene-d12	23.759	264	6224	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	1780	0.662	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.208	152	3756	0.346	ng/ul	0.00
18) Fluoranthene-d10	19.249	212	7896	0.375	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.244	88	4859	1.678	ng/ul#	86
5) Naphthalene	10.651	128	6973	0.347	ng/ul	99
7) 2-Methylnaphthalene	12.279	142	4147	0.338	ng/ul	99
8) 1-Methylnaphthalene	12.493	142	4533	0.335	ng/ul	100
10) Acenaphthylene	14.174	152	5928	0.354	ng/ul	99
11) Acenaphthene	14.521	153	4837	0.346	ng/ul	100
12) Fluorene	15.511	166	5306	0.329	ng/ul	99
14) Pentachlorophenol	16.876	266	762	0.596	ng/ul	98
15) Phenanthrene	17.252	178	8580	0.316	ng/ul	100
16) Anthracene	17.349	178	6970	0.298	ng/ul	100
19) Fluoranthene	19.277	202	9422	0.305	ng/ul	99
20) Pyrene	19.640	202	9764	0.301	ng/ul	96
21) Benzo(a)anthracene	21.401	228	7197	0.288	ng/ul	99
22) Chrysene	21.450	228	9086	0.294	ng/ul	99
24) Benzo(b)fluoranthene	23.043	252	8499	0.286	ng/ul	99
25) Benzo(k)fluoranthene	23.093	252	9168	0.285	ng/ul	99
26) Benzo(a)pyrene	23.654	252	8128	0.290	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.179	276	9054	0.275	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.196	278	7278	0.270	ng/ul	99
29) Benzo(g,h,i)perylene	26.917	276	7965	0.266	ng/ul	98

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

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