

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081022\
 Data File : BN021156.D
 Acq On : 11 Aug 2022 20:24
 Operator : CG/JU
 Sample : PB146935BS
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS935

Quant Time: Aug 11 23:18:37 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 09 23:10:06 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.792	152	2621	0.400	ng/ul	0.01
4) Naphthalene-d8	10.585	136	9182	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.437	164	5255	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.192	188	9821	0.400	ng/ul	0.00
17) Chrysene-d12	21.394	240	6873	0.400	ng/ul	-0.01
23) Perylene-d12	23.727	264	5227	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.197	96	2333	0.696	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.190	152	4869	0.362	ng/ul	0.01
18) Fluoranthene-d10	19.230	212	9307	0.438	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.231	88	6212	1.798	ng/ul#	87
5) Naphthalene	10.634	128	8793	0.312	ng/ul	100
7) 2-Methylnaphthalene	12.262	142	5314	0.311	ng/ul	99
8) 1-Methylnaphthalene	12.476	142	5838	0.341	ng/ul	100
10) Acenaphthylene	14.160	152	7786	0.317	ng/ul	99
11) Acenaphthene	14.502	153	6230	0.307	ng/ul	99
12) Fluorene	15.497	166	6708	0.302	ng/ul	100
14) Pentachlorophenol	16.863	266	754	0.498	ng/ul	100
15) Phenanthrene	17.230	178	10528	0.300	ng/ul	99
16) Anthracene	17.327	178	8636	0.317	ng/ul	99
19) Fluoranthene	19.258	202	11330	0.372	ng/ul	98
20) Pyrene	19.620	202	11753	0.374	ng/ul	97
21) Benzo(a)anthracene	21.380	228	7278	0.325	ng/ul	98
22) Chrysene	21.432	228	9029	0.306	ng/ul	99
24) Benzo(b)fluoranthene	23.019	252	7587	0.314	ng/ul	95
25) Benzo(k)fluoranthene	23.063	252	8084	0.317	ng/ul	98
26) Benzo(a)pyrene	23.624	252	7100	0.322	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	26.135	276	7556	0.287	ng/ul	99
28) Dibenzo(a,h)anthracene	26.151	278	5894	0.283	ng/ul	97
29) Benzo(g,h,i)perylene	26.872	276	6952	0.296	ng/ul	97

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

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