

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081022\
 Data File : BN021127.D
 Acq On : 10 Aug 2022 21:31
 Operator : CG/JU
 Sample : PB146901BS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS901

Quant Time: Aug 10 22:12:53 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	7.793	152	2643	0.400	ng/ul	0.01
4) Naphthalene-d8	10.585	136	9249	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.438	164	5233	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.189	188	9752	0.400	ng/ul	0.00
17) Chrysene-d12	21.395	240	7148	0.400	ng/ul	0.00
23) Perylene-d12	23.725	264	5362	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.198	96	2308	0.683	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.191	152	4816	0.355	ng/ul	0.01
18) Fluoranthene-d10	19.226	212	9407	0.426	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.228	88	6200	1.780	ng/ul#	85
5) Naphthalene	10.635	128	8840	0.311	ng/ul	99
7) 2-Methylnaphthalene	12.263	142	5316	0.308	ng/ul	99
8) 1-Methylnaphthalene	12.477	142	5811	0.337	ng/ul	100
10) Acenaphthylene	14.161	152	7587	0.310	ng/ul	99
11) Acenaphthene	14.503	153	6144	0.304	ng/ul	100
12) Fluorene	15.498	166	6651	0.301	ng/ul	99
14) Pentachlorophenol	16.859	266	718	0.477	ng/ul	100
15) Phenanthrene	17.231	178	10509	0.301	ng/ul	99
16) Anthracene	17.328	178	8516	0.315	ng/ul	99
19) Fluoranthene	19.259	202	11441	0.362	ng/ul	99
20) Pyrene	19.621	202	11708	0.358	ng/ul	98
21) Benzo(a)anthracene	21.380	228	7552	0.324	ng/ul	99
22) Chrysene	21.430	228	9555	0.311	ng/ul	99
24) Benzo(b)fluoranthene	23.017	252	7845	0.316	ng/ul	97
25) Benzo(k)fluoranthene	23.064	252	8393	0.320	ng/ul	98
26) Benzo(a)pyrene	23.622	252	7198	0.318	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.135	276	7516	0.278	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.149	278	5862	0.274	ng/ul	99
29) Benzo(g,h,i)perylene	26.867	276	6912	0.287	ng/ul	98

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

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