

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081022\
 Data File : BN021142.D
 Acq On : 11 Aug 2022 10:56
 Operator : CG/JU
 Sample : PB146918BS
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS918

Quant Time: Aug 11 13:49:23 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	7.789	152	3030	0.400	ng/ul	0.00
4) Naphthalene-d8	10.585	136	10581	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.438	164	6008	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.189	188	11243	0.400	ng/ul	0.00
17) Chrysene-d12	21.398	240	7920	0.400	ng/ul	0.00
23) Perylene-d12	23.727	264	6108	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.198	96	2417	0.624	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.186	152	4970	0.321	ng/ul	0.00
18) Fluoranthene-d10	19.226	212	9470	0.387	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.232	88	6242	1.563	ng/ul#	87
5) Naphthalene	10.629	128	8928	0.275	ng/ul	100
7) 2-Methylnaphthalene	12.263	142	5440	0.276	ng/ul	99
8) 1-Methylnaphthalene	12.477	142	5976	0.303	ng/ul	100
10) Acenaphthylene	14.156	152	7894	0.281	ng/ul	99
11) Acenaphthene	14.498	153	6327	0.272	ng/ul	99
12) Fluorene	15.493	166	6821	0.269	ng/ul	98
14) Pentachlorophenol	16.859	266	744	0.429	ng/ul	97
15) Phenanthrene	17.231	178	10791	0.268	ng/ul	99
16) Anthracene	17.328	178	8696	0.279	ng/ul	99
19) Fluoranthene	19.259	202	11560	0.330	ng/ul	99
20) Pyrene	19.621	202	11928	0.330	ng/ul	96
21) Benzo(a)anthracene	21.383	228	7342	0.284	ng/ul	99
22) Chrysene	21.433	228	9480	0.279	ng/ul	99
24) Benzo(b)fluoranthene	23.020	252	8110	0.287	ng/ul	96
25) Benzo(k)fluoranthene	23.067	252	8232	0.276	ng/ul	98
26) Benzo(a)pyrene	23.625	252	7379	0.286	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.135	276	7746	0.252	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.159	278	5987	0.246	ng/ul	96
29) Benzo(g,h,i)perylene	26.873	276	7081	0.258	ng/ul	97

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

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