

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070821\
 Data File : BN015414.D
 Acq On : 08 Jul 2021 18:39
 Operator : CG/JU
 Sample : PB137471BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS471

Quant Time: Jul 09 01:15:09 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 08 11:17:31 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.711	152	2887	0.400	ng/ul	0.00
4) Naphthalene-d8	10.479	136	9689	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.339	164	4963	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.086	188	10088	0.400	ng/ul	0.00
17) Chrysene-d12	21.281	240	8475	0.400	ng/ul	0.00
23) Perylene-d12	23.520	264	8059	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	2552	0.784	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.074	152	5438	0.358	ng/ul	0.00
18) Fluoranthene-d10	19.117	212	10857	0.390	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.310	88	6536	1.859	ng/ul#	87
5) Naphthalene	10.529	128	10676	0.345	ng/ul	99
7) 2-Methylnaphthalene	12.145	142	6960	0.344	ng/ul	100
8) 1-Methylnaphthalene	12.371	142	6473	0.331	ng/ul	99
10) Acenaphthylene	14.057	152	8273	0.322	ng/ul	99
11) Acenaphthene	14.399	153	6861	0.349	ng/ul	99
12) Fluorene	15.389	166	7624	0.351	ng/ul	99
14) Pentachlorophenol	16.739	266	1427	0.639	ng/ul	99
15) Phenanthrene	17.128	178	13103	0.358	ng/ul	99
16) Anthracene	17.216	178	10927	0.338	ng/ul	100
19) Fluoranthene	19.150	202	14206	0.363	ng/ul	99
20) Pyrene	19.512	202	14384	0.366	ng/ul	99
21) Benzo(a)anthracene	21.266	228	12794	0.408	ng/ul	99
22) Chrysene	21.319	228	13997	0.372	ng/ul	99
24) Benzo(b)fluoranthene	22.850	252	13836	0.367	ng/ul	95
25) Benzo(k)fluoranthene	22.894	252	14485	0.341	ng/ul	98
26) Benzo(a)pyrene	23.423	252	13117	0.415	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	25.770	276	15546	0.407	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.783	278	12397	0.414	ng/ul	99
29) Benzo(g,h,i)perylene	26.457	276	14223	0.402	ng/ul	98

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

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