

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062121\
 Data File : BN014961.D
 Acq On : 21 Jun 2021 19:39
 Operator : CG/JU
 Sample : PB137168BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS168

Quant Time: Jun 22 02:16:03 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 : Mon Jun 21 14:33:43 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.743	152	3385	0.400	ng/ul	0.00
4) Naphthalene-d8	10.516	136	11395	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.366	164	5632	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.114	188	11138	0.400	ng/ul	0.00
17) Chrysene-d12	21.312	240	8487	0.400	ng/ul	0.00
23) Perylene-d12	23.569	264	7000	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.300	96	2907	0.762	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.105	152	6710	0.375	ng/ul	0.00
18) Fluoranthene-d10	19.149	212	12478	0.447	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.330	88	7706	1.869	ng/ul#	86
5) Naphthalene	10.565	128	13491	0.371	ng/ul	100
7) 2-Methylnaphthalene	12.182	142	8617	0.362	ng/ul	98
8) 1-Methylnaphthalene	12.402	142	8066	0.351	ng/ul	99
10) Acenaphthylene	14.084	152	10247	0.352	ng/ul	100
11) Acenaphthene	14.431	153	8232	0.369	ng/ul	99
12) Fluorene	15.421	166	8974	0.364	ng/ul	99
14) Pentachlorophenol	16.760	266	1333	0.541	ng/ul	100
15) Phenanthrene	17.157	178	15224	0.377	ng/ul	99
16) Anthracene	17.249	178	12680	0.356	ng/ul	99
19) Fluoranthene	19.177	202	16369	0.417	ng/ul	97
20) Pyrene	19.540	202	16381	0.416	ng/ul	97
21) Benzo(a)anthracene	21.298	228	12269	0.390	ng/ul	99
22) Chrysene	21.350	228	15363	0.408	ng/ul	100
24) Benzo(b)fluoranthene	22.893	252	12480	0.382	ng/ul	98
25) Benzo(k)fluoranthene	22.940	252	14651	0.397	ng/ul	97
26) Benzo(a)pyrene	23.469	252	10267	0.374	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.849	276	13205	0.398	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.866	278	10368	0.399	ng/ul	98
29) Benzo(g,h,i)perylene	26.534	276	12109	0.394	ng/ul	96

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

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