

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082824\
 Data File : BN033652.D
 Acq On : 28 Aug 2024 18:59
 Operator : MA/JU
 Sample : PB162963BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS963

Quant Time: Aug 28 23:10:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 23:48:09 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.536	152	5127	0.400	ng/ul	0.00
4) Naphthalene-d8	10.294	136	13766	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.174	164	6841	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.927	188	14285	0.400	ng/ul	0.00
17) Chrysene-d12	21.138	240	11403	0.400	ng/ul	0.00
23) Perylene-d12	23.301	264	12379	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	4147	0.787	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.894	152	7911	0.401	ng/ul	0.00
18) Fluoranthene-d10	18.966	212	14566	0.428	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.177	88	11461	1.960	ng/ul#	79
5) Naphthalene	10.343	128	14154	0.384	ng/ul	99
7) 2-Methylnaphthalene	11.971	142	8908	0.382	ng/ul	99
8) 1-Methylnaphthalene	12.191	142	10569	0.441	ng/ul	99
10) Acenaphthylene	13.892	152	11461	0.369	ng/ul	99
11) Acenaphthene	14.239	153	9216	0.411	ng/ul	99
12) Fluorene	15.234	166	9906	0.395	ng/ul	99
14) Pentachlorophenol	16.585	266	3007	0.669	ng/ul	99
15) Phenanthrene	16.965	178	15492	0.368	ng/ul	100
16) Anthracene	17.058	178	13951	0.368	ng/ul	99
19) Fluoranthene	18.994	202	17399	0.387	ng/ul	99
20) Pyrene	19.361	202	18217	0.381	ng/ul	99
21) Benzo(a)anthracene	21.121	228	16891	0.376	ng/ul	99
22) Chrysene	21.174	228	17478	0.385	ng/ul	99
24) Benzo(b)fluoranthene	22.658	252	17709	0.369	ng/ul	100
25) Benzo(k)fluoranthene	22.702	252	18400	0.383	ng/ul	100
26) Benzo(a)pyrene	23.208	252	16011	0.415	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.448	276	21759	0.375	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.465	278	16977	0.377	ng/ul	99
29) Benzo(g,h,i)perylene	26.102	276	17275	0.381	ng/ul	99

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

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