

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111924\
 Data File : BN035193.D
 Acq On : 20 Nov 2024 11:31
 Operator : RC/JU
 Sample : PB164843BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS843

Quant Time: Nov 20 11:57:50 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 18 00:44:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.541	152	3066	0.400	ng/ul	0.00
4) Naphthalene-d8	10.301	136	7386	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.180	164	5000	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.937	188	10307	0.400	ng/ul	0.00
17) Chrysene-d12	21.148	240	7421	0.400	ng/ul	0.00
23) Perylene-d12	23.320	264	7779	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.136	96	1829	0.653	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.901	152	4041	0.373	ng/ul	-0.01
18) Fluoranthene-d10	18.977	212	8520	0.416	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.170	88	4909	1.614	ng/ul#	95
5) Naphthalene	10.350	128	6574	0.353	ng/ul	99
7) 2-Methylnaphthalene	11.978	142	4618	0.358	ng/ul	100
8) 1-Methylnaphthalene	12.198	142	4648	0.357	ng/ul	99
10) Acenaphthylene	13.898	152	6845	0.345	ng/ul	100
11) Acenaphthene	14.245	153	5299	0.349	ng/ul	98
12) Fluorene	15.240	166	6099	0.342	ng/ul	100
14) Pentachlorophenol	16.595	266	1367	0.569	ng/ul	97
15) Phenanthrene	16.979	178	9386	0.344	ng/ul	100
16) Anthracene	17.068	178	8500	0.336	ng/ul	100
19) Fluoranthene	19.005	202	10157	0.393	ng/ul	100
20) Pyrene	19.367	202	10364	0.385	ng/ul	98
21) Benzo(a)anthracene	21.134	228	9160	0.356	ng/ul	100
22) Chrysene	21.183	228	9629	0.372	ng/ul	99
24) Benzo(b)fluoranthene	22.674	252	9592	0.373	ng/ul	99
25) Benzo(k)fluoranthene	22.718	252	9799	0.370	ng/ul	100
26) Benzo(a)pyrene	23.226	252	8589	0.366	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.477	276	11740	0.370	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.493	278	9162	0.368	ng/ul	99
29) Benzo(g,h,i)perylene	26.127	276	9302	0.372	ng/ul	99

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

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