

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042023\
 Data File : BN025044.D
 Acq On : 20 Apr 2023 20:25
 Operator : CG/JU
 Sample : PB152218BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS218

Quant Time: Apr 21 04:27:17 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 21 04:13:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	5900	0.400	ng/u1	0.00
4) Naphthalene-d8	10.743	136	18522	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.575	164	10315	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.321	188	22052	0.400	ng/u1	0.00
17) Chrysene-d12	21.514	240	15299	0.400	ng/u1	0.00
23) Perylene-d12	23.961	264	11788	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	5010	0.820	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.332	152	9670	0.407	ng/u1	0.00
18) Fluoranthene-d10	19.349	212	20396	0.506	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.339	88	12661	1.924	ng/u1#	87
5) Naphthalene	10.792	128	18507	0.390	ng/u1	99
7) 2-Methylnaphthalene	12.403	142	11373	0.391	ng/u1	99
8) 1-Methylnaphthalene	12.623	142	12156	0.392	ng/u1	100
10) Acenaphthylene	14.297	152	15096	0.393	ng/u1	100
11) Acenaphthene	14.635	153	13186	0.407	ng/u1	99
12) Fluorene	15.625	166	14548	0.402	ng/u1	99
14) Pentachlorophenol	16.975	266	2800	0.647	ng/u1	99
15) Phenanthrene	17.364	178	24185	0.389	ng/u1	100
16) Anthracene	17.461	178	19137	0.376	ng/u1	99
19) Fluoranthene	19.382	202	26683	0.482	ng/u1	99
20) Pyrene	19.744	202	27049	0.476	ng/u1	100
21) Benzo(a)anthracene	21.500	228	19233	0.417	ng/u1	99
22) Chrysene	21.552	228	22303	0.443	ng/u1	100
24) Benzo(b)fluoranthene	23.212	252	20152	0.421	ng/u1	99
25) Benzo(k)fluoranthene	23.262	252	19499	0.424	ng/u1	96
26) Benzo(a)pyrene	23.859	252	15644	0.395	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	26.514	276	18648	0.409	ng/u1	99
28) Dibenzo(a,h)anthracene	26.531	278	14293	0.401	ng/u1	98
29) Benzo(g,h,i)perylene	27.292	276	17185	0.443	ng/u1	99

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

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