

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040424\
 Data File : BM044918.D
 Acq On : 04 Apr 2024 17:12
 Operator : MA/JU
 Sample : PB159946BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS946

Quant Time: Apr 04 18:12:46 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM040424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 03 23:34:03 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.652	152	2305	0.400	ng/ul	0.00
4) Naphthalene-d8	10.430	136	7759	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.298	164	4065	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.061	188	8054	0.400	ng/ul	0.00
17) Chrysene-d12	21.268	240	7473	0.400	ng/ul	# 0.00
23) Perylene-d12	23.498	264	9290	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.226	96	1871	0.618	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.030	152	3904	0.371	ng/ul	-0.02
18) Fluoranthene-d10	19.099	212	7917	0.419	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.255	88	5215	1.797	ng/ul#	84
5) Naphthalene	10.479	128	7715	0.375	ng/ul	99
7) 2-Methylnaphthalene	12.107	142	4816	0.392	ng/ul	95
8) 1-Methylnaphthalene	12.321	142	5033	0.375	ng/ul	99
10) Acenaphthylene	14.021	152	8205	0.405	ng/ul	99
11) Acenaphthene	14.363	153	5578	0.384	ng/ul	99
12) Fluorene	15.362	166	5947	0.365	ng/ul	100
14) Pentachlorophenol	16.719	266	1780	0.943	ng/ul	96
15) Phenanthrene	17.103	178	9136	0.403	ng/ul	99
16) Anthracene	17.196	178	9326	0.418	ng/ul	100
19) Fluoranthene	19.127	202	11669	0.437	ng/ul	99
20) Pyrene	19.495	202	12462	0.442	ng/ul	100
21) Benzo(a)anthracene	21.254	228	11274	0.433	ng/ul	93
22) Chrysene	21.303	228	12318	0.379	ng/ul#	93
24) Benzo(b)fluoranthene	22.829	252	12513	0.373	ng/ul#	73
25) Benzo(k)fluoranthene	22.873	252	13347	0.355	ng/ul#	74
26) Benzo(a)pyrene	23.399	252	12402	0.387	ng/ul#	71
27) Indeno(1,2,3-cd)pyrene	25.745	276	16119	0.367	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.765	278	13012	0.374	ng/ul#	75
29) Benzo(g,h,i)perylene	26.429	276	13405	0.354	ng/ul#	82

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

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