

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042623\
 Data File : BM039681.D
 Acq On : 26 Apr 2023 16:45
 Operator : CG/JU
 Sample : PB152372BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS372

Quant Time: Apr 27 01:03:25 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM040623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 26 19:06:24 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	2127	0.400	ng/ul	0.00
4) Naphthalene-d8	10.639	136	11028	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.483	164	6044	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.250	188	9079	0.400	ng/ul	0.00
17) Chrysene-d12	21.442	240	7610	0.400	ng/ul	# 0.00
23) Perylene-d12	23.812	264	6386	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.214	96	3551	0.902	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	5526	0.433	ng/ul	0.02
18) Fluoranthene-d10	19.280	212	9440	0.398	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.247	88	8975	2.129	ng/ul#	78
5) Naphthalene	10.694	128	10960	0.418	ng/ul	98
7) 2-Methylnaphthalene	12.327	142	6022	0.409	ng/ul	95
8) 1-Methylnaphthalene	12.525	142	6502	0.390	ng/ul	98
10) Acenaphthylene	14.210	152	10363	0.442	ng/ul	98
11) Acenaphthene	14.548	153	8061	0.436	ng/ul	99
12) Fluorene	15.552	166	8173	0.444	ng/ul	98
14) Pentachlorophenol	16.951	266	1285	0.594	ng/ul#	81
15) Phenanthrene	17.297	178	10829	0.430	ng/ul	97
16) Anthracene	17.394	178	8048	0.420	ng/ul	98
19) Fluoranthene	19.308	202	12370	0.393	ng/ul	98
20) Pyrene	19.671	202	14094	0.390	ng/ul	99
21) Benzo(a)anthracene	21.430	228	7616	0.434	ng/ul	98
22) Chrysene	21.477	228	10706	0.498	ng/ul	100
24) Benzo(b)fluoranthene	23.087	252	9926	0.490	ng/ul	95
25) Benzo(k)fluoranthene	23.137	252	9735	0.504	ng/ul	95
26) Benzo(a)pyrene	23.710	252	9467	0.464	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.286	276	13062	0.508	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.300	278	9950	0.553	ng/ul	96
29) Benzo(g,h,i)perylene	27.024	276	11895	0.482	ng/ul	99

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

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