

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071123\
 Data File : BM040814.D
 Acq On : 11 Jul 2023 15:59
 Operator : MA/JU
 Sample : PB154053BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS053

Quant Time: Jul 12 04:34:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 11 03:54:24 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	2689	0.400	ng/ul	0.00
4) Naphthalene-d8	10.691	136	10484	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.532	164	4636	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.291	188	7472	0.400	ng/ul	0.00
17) Chrysene-d12	21.483	240	5352	0.400	ng/ul	0.00
23) Perylene-d12	23.874	264	4376	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.297	96	3311	0.784	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.286	152	5096	0.347	ng/ul	0.00
18) Fluoranthene-d10	19.320	212	6232	0.363	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.330	88	7394	1.720	ng/ul	90
5) Naphthalene	10.741	128	9886	0.342	ng/ul	97
7) 2-Methylnaphthalene	12.357	142	5580	0.329	ng/ul	97
8) 1-Methylnaphthalene	12.572	142	5540	0.326	ng/ul	98
10) Acenaphthylene	14.254	152	6992	0.338	ng/ul	94
11) Acenaphthene	14.597	153	5837	0.334	ng/ul	99
12) Fluorene	15.591	166	6031	0.323	ng/ul	99
14) Pentachlorophenol	16.949	266	422	0.293	ng/ul	90
15) Phenanthrene	17.333	178	7599	0.327	ng/ul#	95
16) Anthracene	17.426	178	7442	0.382	ng/ul#	93
19) Fluoranthene	19.348	202	7872	0.346	ng/ul	96
20) Pyrene	19.710	202	8662	0.355	ng/ul#	95
21) Benzo(a)anthracene	21.466	228	6367	0.352	ng/ul	96
22) Chrysene	21.518	228	7420	0.359	ng/ul	98
24) Benzo(b)fluoranthene	23.146	252	6260	0.370	ng/ul	82
25) Benzo(k)fluoranthene	23.190	252	6812	0.408	ng/ul#	87
26) Benzo(a)pyrene	23.769	252	5944	0.390	ng/ul#	80
27) Indeno(1,2,3-cd)pyrene	26.357	276	8453	0.398	ng/ul#	44
28) Dibenzo(a,h)anthracene	26.374	278	6149	0.370	ng/ul#	89
29) Benzo(g,h,i)perylene	27.112	276	6578	0.353	ng/ul	94

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

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