

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM063023\
 Data File : BM040476.D
 Acq On : 30 Jun 2023 17:44
 Operator : MA/JU
 Sample : PB153575BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS575

Quant Time: Jul 01 03:04:13 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 : Sat Jul 01 03:00:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	5310	0.400	ng/ul	0.00
4) Naphthalene-d8	10.749	136	20633	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.580	164	10562	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.331	188	20488	0.400	ng/ul	0.00
17) Chrysene-d12	21.509	240	13956	0.400	ng/ul	0.00
23) Perylene-d12	23.924	264	12547	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.325	96	3622	0.434	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.338	152	7894	0.273	ng/ul	0.00
18) Fluoranthene-d10	19.355	212	12793	0.286	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.363	88	9897	1.166	ng/ul	90
5) Naphthalene	10.799	128	15397	0.271	ng/ul	98
7) 2-Methylnaphthalene	12.410	142	9154	0.274	ng/ul	99
8) 1-Methylnaphthalene	12.630	142	9156	0.274	ng/ul	99
10) Acenaphthylene	14.303	152	12678	0.269	ng/ul	99
11) Acenaphthene	14.645	153	10201	0.256	ng/ul	99
12) Fluorene	15.630	166	10911	0.257	ng/ul	100
14) Pentachlorophenol	16.985	266	581	0.147	ng/ul	92
15) Phenanthrene	17.373	178	15984	0.251	ng/ul	99
16) Anthracene	17.462	178	13944	0.261	ng/ul	98
19) Fluoranthene	19.387	202	16792	0.283	ng/ul	97
20) Pyrene	19.750	202	17920	0.282	ng/ul	99
21) Benzo(a)anthracene	21.492	228	13987	0.296	ng/ul	99
22) Chrysene	21.547	228	14446	0.268	ng/ul	99
24) Benzo(b)fluoranthene	23.184	252	13330	0.275	ng/ul	91
25) Benzo(k)fluoranthene	23.234	252	13137	0.275	ng/ul#	90
26) Benzo(a)pyrene	23.815	252	11454	0.262	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	26.427	276	14919	0.245	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.444	278	11894	0.249	ng/ul	94
29) Benzo(g,h,i)perylene	27.192	276	12553	0.235	ng/ul	98

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

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