

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071124\
 Data File : BM046536.D
 Acq On : 12 Jul 2024 03:05
 Operator : MA/JU
 Sample : PB161755BS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS755

Quant Time: Jul 12 05:12:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 05:09:19 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.510	152	2913	0.400	ng/ul	0.00
4) Naphthalene-d8	10.271	136	7282	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.155	164	4577	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.905	188	9849	0.400	ng/ul	0.00
17) Chrysene-d12	21.111	240	7566	0.400	ng/ul	0.00
23) Perylene-d12	23.250	264	8707	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	1327	0.554	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.871	152	3608	0.307	ng/ul	0.00
18) Fluoranthene-d10	18.942	212	7711	0.335	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.151	88	3492	1.367	ng/ul#	69
5) Naphthalene	10.321	128	5465	0.295	ng/ul	97
7) 2-Methylnaphthalene	11.948	142	3843	0.295	ng/ul	99
8) 1-Methylnaphthalene	12.168	142	3874	0.292	ng/ul	97
10) Acenaphthylene	13.868	152	5991	0.288	ng/ul	98
11) Acenaphthene	14.215	153	4318	0.309	ng/ul	98
12) Fluorene	15.210	166	5064	0.293	ng/ul	98
14) Pentachlorophenol	16.559	266	2011	0.403	ng/ul	97
15) Phenanthrene	16.947	178	8146	0.290	ng/ul	99
16) Anthracene	17.036	178	7806	0.285	ng/ul	99
19) Fluoranthene	18.974	202	9902	0.319	ng/ul	99
20) Pyrene	19.337	202	10269	0.305	ng/ul	99
21) Benzo(a)anthracene	21.093	228	9707	0.290	ng/ul	99
22) Chrysene	21.146	228	9587	0.296	ng/ul	99
24) Benzo(b)fluoranthene	22.616	252	10433	0.300	ng/ul	97
25) Benzo(k)fluoranthene	22.659	252	10040	0.294	ng/ul	97
26) Benzo(a)pyrene	23.159	252	9224	0.322	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.373	276	12568	0.289	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.393	278	9861	0.286	ng/ul	99
29) Benzo(g,h,i)perylene	26.017	276	9926	0.286	ng/ul	99

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

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