

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072524\
 Data File : BM046793.D
 Acq On : 25 Jul 2024 10:35
 Operator : MA/JU
 Sample : P3263-25
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4DQ6

Quant Time: Jul 25 11:11:46 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM072424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 24 15:10:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.472	152	2158	0.400	ng/ul	0.00
4) Naphthalene-d8	10.233	136	5611	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.118	164	3773	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.871	188	8704	0.400	ng/ul	0.00
17) Chrysene-d12	21.082	240	7095	0.400	ng/ul	0.00
23) Perylene-d12	23.209	264	7164	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.101	96	574	0.380	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.827	152	2298	0.258	ng/ul	0.00
18) Fluoranthene-d10	18.909	212	6850	0.296	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.277	128	194406	13.698	ng/ul	96
10) Acenaphthylene	13.831	152	222587	12.881	ng/ul	97
11) Acenaphthene	14.178	153	137499	11.723	ng/ul	100
12) Fluorene	15.178	166	95042	6.483	ng/ul	96
15) Phenanthrene	16.913	178	156120	6.240	ng/ul	97
16) Anthracene	17.002	178	113447	4.871	ng/ul	99
19) Fluoranthene	18.942	202	223741	7.180	ng/ul#	97
20) Pyrene	19.304	202	274781	8.525	ng/ul	97
21) Benzo(a)anthracene	21.067	228	70473	2.384	ng/ul	99
22) Chrysene	21.120	228	124464	4.144	ng/ul	98
24) Benzo(b)fluoranthene	22.580	252	145002	4.610	ng/ul	90
25) Benzo(k)fluoranthene	22.621	252	121943	3.847	ng/ul#	87
26) Benzo(a)pyrene	23.118	252	144240	6.388	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	25.313	276	166798	4.800	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.329	278	102264	4.020	ng/ul#	87
29) Benzo(g,h,i)perylene	25.953	276	119777	3.914	ng/ul	96

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

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