

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072524\
 Data File : BM046794.D
 Acq On : 25 Jul 2024 11:40
 Operator : MA/JU
 Sample : P3263-25DL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4DQ6DL

Quant Time: Jul 25 12:11:52 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	3341	0.400	ng/ul	0.00
4) Naphthalene-d8	10.227	136	8951	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.118	164	5713	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.871	188	12110	0.400	ng/ul	0.00
17) Chrysene-d12	21.082	240	7267	0.400	ng/ul	0.00
23) Perylene-d12	23.209	264	6943	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.101	96	143	0.061	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.827	152	361	0.025	ng/ul	0.00
18) Fluoranthene-d10	18.914	212	879	0.037	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.277	128	31257	1.381	ng/ul	97
10) Acenaphthylene	13.831	152	33249	1.271	ng/ul	98
11) Acenaphthene	14.178	153	20761	1.169	ng/ul	99
12) Fluorene	15.178	166	13801	0.622	ng/ul	97
15) Phenanthrene	16.913	178	21977	0.631	ng/ul	98
16) Anthracene	17.006	178	14795	0.457	ng/ul	100
19) Fluoranthene	18.942	202	27694	0.868	ng/ul	99
20) Pyrene	19.309	202	33023	1.000	ng/ul	97
21) Benzo(a)anthracene	21.067	228	7202	0.238	ng/ul	99
22) Chrysene	21.119	228	13063	0.425	ng/ul	99
24) Benzo(b)fluoranthene	22.580	252	13906	0.456	ng/ul	98
25) Benzo(k)fluoranthene	22.621	252	12303	0.401	ng/ul	98
26) Benzo(a)pyrene	23.115	252	13989	0.639	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.313	276	20052	0.595	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.326	278	12354	0.501	ng/ul	99
29) Benzo(g,h,i)perylene	25.950	276	11593	0.391	ng/ul	98

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

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