

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072524\
 Data File : BM046818.D
 Acq On : 26 Jul 2024 02:18
 Operator : MA/JU
 Sample : PB162156BS
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS156

Quant Time: Jul 26 02:51:46 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	3238	0.400	ng/ul	0.00
4) Naphthalene-d8	10.226	136	9439	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.117	164	5942	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.874	188	11760	0.400	ng/ul	0.00
17) Chrysene-d12	21.082	240	6608	0.400	ng/ul	# 0.00
23) Perylene-d12	23.212	264	7007	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.101	96	1596	0.704	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.826	152	6571	0.438	ng/ul	0.00
18) Fluoranthene-d10	18.913	212	11347	0.527	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.130	88	4311	1.662	ng/ul#	83
5) Naphthalene	10.275	128	10267	0.430	ng/ul	99
7) 2-Methylnaphthalene	11.903	142	7419	0.452	ng/ul	99
8) 1-Methylnaphthalene	12.129	142	7489	0.441	ng/ul	99
10) Acenaphthylene	13.830	152	12108	0.445	ng/ul	99
11) Acenaphthene	14.182	153	8483	0.459	ng/ul	98
12) Fluorene	15.176	166	9699	0.420	ng/ul	97
14) Pentachlorophenol	16.532	266	3665	0.688	ng/ul	99
15) Phenanthrene	16.912	178	14336	0.424	ng/ul	98
16) Anthracene	17.005	178	13625	0.433	ng/ul	99
19) Fluoranthene	18.945	202	15327	0.528	ng/ul	99
20) Pyrene	19.308	202	15610	0.520	ng/ul	98
21) Benzo(a)anthracene	21.067	228	12360	0.449	ng/ul	99
22) Chrysene	21.117	228	11895	0.425	ng/ul	98
24) Benzo(b)fluoranthene	22.578	252	11861	0.386	ng/ul	90
25) Benzo(k)fluoranthene	22.622	252	11812	0.381	ng/ul#	94
26) Benzo(a)pyrene	23.116	252	10756	0.487	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	25.316	276	14199	0.418	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.333	278	10896	0.438	ng/ul#	94
29) Benzo(g,h,i)perylene	25.957	276	11215	0.375	ng/ul#	94

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

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