

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072924\
 Data File : BM046918.D
 Acq On : 30 Jul 2024 12:52
 Operator : MA/JU
 Sample : PB162178BS
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS178

Quant Time: Jul 30 13:38:20 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 : Tue Jul 30 04:18:24 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.451	152	2545	0.400	ng/ul	0.00
4) Naphthalene-d8	10.205	136	14364	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.100	164	4177	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.854	188	7608	0.400	ng/ul	0.00
17) Chrysene-d12	21.073	240	4948	0.400	ng/ul	0.00
23) Perylene-d12	23.197	264	5738	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.084	96	1273	0.714	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.811	152	4882	0.214	ng/ul	0.00
18) Fluoranthene-d10	18.900	212	7277	0.452	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.117	88	3461	1.697	ng/ul#	85
5) Naphthalene	10.255	128	14798	0.407	ng/ul	98
7) 2-Methylnaphthalene	11.882	142	5190	0.208	ng/ul	100
8) 1-Methylnaphthalene	12.108	142	5184	0.201	ng/ul	99
10) Acenaphthylene	13.813	152	7870	0.411	ng/ul	99
11) Acenaphthene	14.160	153	5546	0.427	ng/ul	97
12) Fluorene	15.159	166	6237	0.384	ng/ul	95
14) Pentachlorophenol	16.516	266	2235	0.649	ng/ul	98
15) Phenanthrene	16.896	178	8560	0.391	ng/ul	100
16) Anthracene	16.989	178	8007	0.393	ng/ul	99
19) Fluoranthene	18.928	202	8979	0.413	ng/ul	99
20) Pyrene	19.295	202	9249	0.411	ng/ul	98
21) Benzo(a)anthracene	21.058	228	8011	0.389	ng/ul	100
22) Chrysene	21.111	228	8037	0.384	ng/ul	98
24) Benzo(b)fluoranthene	22.569	252	8454	0.336	ng/ul	97
25) Benzo(k)fluoranthene	22.610	252	9010	0.355	ng/ul#	97
26) Benzo(a)pyrene	23.104	252	7789	0.431	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.292	276	11106	0.399	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.309	278	8565	0.420	ng/ul	99
29) Benzo(g,h,i)perylene	25.930	276	8947	0.365	ng/ul	99

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

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