

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082923\
 Data File : BM041576.D
 Acq On : 30 Aug 2023 12:42
 Operator : MA/JU
 Sample : PB155079BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS079

Quant Time: Aug 31 02:16:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM082923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 30 01:50:06 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.114	152	6356	0.400	ng/ul	0.00
4) Naphthalene-d8	10.944	136	14265	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.754	164	5523	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.506	188	10080	0.400	ng/ul	0.00
17) Chrysene-d12	21.688	240	3538	0.400	ng/ul	0.00
23) Perylene-d12	24.249	264	3654	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.452	96	6035	0.643	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.522	152	6015	0.350	ng/ul	0.00
18) Fluoranthene-d10	19.524	212	5996	0.344	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.490	88	14903	1.489	ng/ul#	85
5) Naphthalene	10.999	128	14654	0.321	ng/ul	100
7) 2-Methylnaphthalene	12.594	142	8109	0.313	ng/ul	100
8) 1-Methylnaphthalene	12.808	142	8471	0.324	ng/ul	99
10) Acenaphthylene	14.481	152	12157	0.316	ng/ul	100
11) Acenaphthene	14.814	153	8414	0.317	ng/ul	98
12) Fluorene	15.800	166	9032	0.312	ng/ul	97
14) Pentachlorophenol	17.126	266	2061	0.532	ng/ul	98
15) Phenanthrene	17.544	178	13706	0.311	ng/ul	99
16) Anthracene	17.637	178	11801	0.309	ng/ul	99
19) Fluoranthene	19.552	202	13826	0.298	ng/ul	97
20) Pyrene	19.919	202	13686	0.293	ng/ul	97
21) Benzo(a)anthracene	21.670	228	9228	0.292	ng/ul	99
22) Chrysene	21.726	228	10143	0.299	ng/ul	99
24) Benzo(b)fluoranthene	23.453	252	10118	0.298	ng/ul	97
25) Benzo(k)fluoranthene	23.509	252	9870	0.306	ng/ul	97
26) Benzo(a)pyrene	24.135	252	8393	0.278	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.958	276	11454	0.284	ng/ul#	37
28) Dibenzo(a,h)anthracene	26.995	278	7741	0.285	ng/ul	98
29) Benzo(g,h,i)perylene	27.803	276	10211	0.301	ng/ul	99

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

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