

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082423\
 Data File : BM041486.D
 Acq On : 25 Aug 2023 04:22
 Operator : MA/JU
 Sample : PB155029BS
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS029

Quant Time: Aug 25 05:01:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM082423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 24 18:34:12 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.114	152	12478	0.400	ng/ul	0.00
4) Naphthalene-d8	10.944	136	33258	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.754	164	15551	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.506	188	30865	0.400	ng/ul	0.00
17) Chrysene-d12	21.697	240	9427	0.400	ng/ul	0.00
23) Perylene-d12	24.266	264	7995	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.448	96	9404	0.595	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.522	152	14050	0.320	ng/ul	0.00
18) Fluoranthene-d10	19.529	212	16223	0.390	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.486	88	22925	1.411	ng/ul#	91
5) Naphthalene	10.994	128	29345	0.296	ng/ul	100
7) 2-Methylnaphthalene	12.594	142	17590	0.295	ng/ul	100
8) 1-Methylnaphthalene	12.814	142	18804	0.311	ng/ul	100
10) Acenaphthylene	14.481	152	26970	0.289	ng/ul	100
11) Acenaphthene	14.819	153	19703	0.303	ng/ul	99
12) Fluorene	15.804	166	22087	0.304	ng/ul	98
14) Pentachlorophenol	17.126	266	4000	0.687	ng/ul	97
15) Phenanthrene	17.548	178	35269	0.311	ng/ul	99
16) Anthracene	17.641	178	30157	0.292	ng/ul	99
19) Fluoranthene	19.562	202	34816	0.402	ng/ul	99
20) Pyrene	19.924	202	34810	0.399	ng/ul	97
21) Benzo(a)anthracene	21.679	228	21432	0.349	ng/ul	98
22) Chrysene	21.737	228	23836	0.360	ng/ul	100
24) Benzo(b)fluoranthene	23.471	252	20440	0.373	ng/ul	98
25) Benzo(k)fluoranthene	23.524	252	19997	0.365	ng/ul	98
26) Benzo(a)pyrene	24.152	252	16964	0.377	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.985	276	19154	0.341	ng/ul#	97
28) Dibenzo(a,h)anthracene	27.021	278	13740	0.319	ng/ul	98
29) Benzo(g,h,i)perylene	27.826	276	16433	0.332	ng/ul	99

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

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