

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082423\
 Data File : BM041466.D
 Acq On : 24 Aug 2023 15:31
 Operator : MA/JU
 Sample : PB155026BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS026

Quant Time: Aug 24 18:36:41 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	8.118	152	10348	0.400	ng/ul	0.00
4) Naphthalene-d8	10.950	136	28520	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.759	164	13012	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.515	188	24803	0.400	ng/ul	0.00
17) Chrysene-d12	21.705	240	7423	0.400	ng/ul	0.00
23) Perylene-d12	24.281	264	7212	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.448	96	8074	0.616	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.528	152	12487	0.332	ng/ul	0.00
18) Fluoranthene-d10	19.534	212	13123	0.401	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.486	88	18992	1.409	ng/ul#	92
5) Naphthalene	10.999	128	24531	0.289	ng/ul	100
7) 2-Methylnaphthalene	12.599	142	14493	0.284	ng/ul	100
8) 1-Methylnaphthalene	12.819	142	15433	0.297	ng/ul	100
10) Acenaphthylene	14.486	152	21811	0.279	ng/ul	100
11) Acenaphthene	14.823	153	15586	0.287	ng/ul	97
12) Fluorene	15.804	166	17145	0.282	ng/ul	100
14) Pentachlorophenol	17.130	266	2168	0.464	ng/ul	98
15) Phenanthrene	17.557	178	26169	0.287	ng/ul	100
16) Anthracene	17.645	178	22593	0.273	ng/ul	100
19) Fluoranthene	19.566	202	24786	0.364	ng/ul	100
20) Pyrene	19.929	202	24875	0.362	ng/ul	96
21) Benzo(a)anthracene	21.688	228	15889	0.329	ng/ul	99
22) Chrysene	21.743	228	18174	0.348	ng/ul	99
24) Benzo(b)fluoranthene	23.483	252	17081	0.346	ng/ul	96
25) Benzo(k)fluoranthene	23.535	252	17260	0.349	ng/ul	99
26) Benzo(a)pyrene	24.164	252	14404	0.355	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	27.008	276	16780	0.331	ng/ul#	95
28) Dibenzo(a,h)anthracene	27.042	278	13147	0.339	ng/ul	98
29) Benzo(g,h,i)perylene	27.850	276	14850	0.333	ng/ul	97

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

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