

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032224\
 Data File : BM044728.D
 Acq On : 22 Mar 2024 09:08
 Operator : MA/JU
 Sample : PB159722BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS722

Quant Time: Mar 22 10:44:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM031124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 20 10:51:03 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.694	152	3862	0.400	ng/ul	0.00
4) Naphthalene-d8	10.474	136	12103	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.335	164	5902	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.098	188	11444	0.400	ng/ul	0.00
17) Chrysene-d12	21.293	240	9078	0.400	ng/ul	0.00
23) Perylene-d12	23.534	264	10839	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.260	96	3933	0.777	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.074	152	6555	0.377	ng/ul	0.00
18) Fluoranthene-d10	19.127	212	11083	0.348	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.289	88	10940	2.115	ng/ul#	65
5) Naphthalene	10.523	128	13208	0.388	ng/ul	98
7) 2-Methylnaphthalene	12.145	142	7771	0.361	ng/ul	98
8) 1-Methylnaphthalene	12.365	142	8022	0.369	ng/ul	100
10) Acenaphthylene	14.057	152	12213	0.407	ng/ul	98
11) Acenaphthene	14.399	153	8558	0.395	ng/ul	99
12) Fluorene	15.399	166	9161	0.370	ng/ul	99
14) Pentachlorophenol	16.752	266	2706	0.980	ng/ul	99
15) Phenanthrene	17.136	178	12982	0.378	ng/ul	98
16) Anthracene	17.233	178	14073	0.436	ng/ul	97
19) Fluoranthene	19.159	202	15028	0.338	ng/ul	96
20) Pyrene	19.522	202	15836	0.348	ng/ul	96
21) Benzo(a)anthracene	21.278	228	13662	0.395	ng/ul	98
22) Chrysene	21.331	228	15715	0.405	ng/ul	99
24) Benzo(b)fluoranthene	22.859	252	15306	0.370	ng/ul	91
25) Benzo(k)fluoranthene	22.903	252	17629	0.393	ng/ul#	92
26) Benzo(a)pyrene	23.435	252	15338	0.408	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	25.793	276	20021	0.463	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.820	278	15798	0.474	ng/ul	96
29) Benzo(g,h,i)perylene	26.481	276	17090	0.450	ng/ul	97

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

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