

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032024\
 Data File : BM044678.D
 Acq On : 20 Mar 2024 11:47
 Operator : MA/JU
 Sample : PB159669BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS669

Quant Time: Mar 20 12:15:44 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	7.702	152	4011	0.400	ng/ul	0.00
4) Naphthalene-d8	10.484	136	12157	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.344	164	5824	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.106	188	11409	0.400	ng/ul	0.00
17) Chrysene-d12	21.304	240	8752	0.400	ng/ul	0.00
23) Perylene-d12	23.552	264	10335	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	3636	0.692	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.079	152	5939	0.340	ng/ul	0.00
18) Fluoranthene-d10	19.140	212	10306	0.335	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.293	88	10629	1.978	ng/ul#	67
5) Naphthalene	10.534	128	11790	0.345	ng/ul	99
7) 2-Methylnaphthalene	12.156	142	7130	0.330	ng/ul	99
8) 1-Methylnaphthalene	12.371	142	7224	0.331	ng/ul	99
10) Acenaphthylene	14.066	152	10710	0.362	ng/ul	100
11) Acenaphthene	14.408	153	7722	0.361	ng/ul	100
12) Fluorene	15.408	166	8434	0.345	ng/ul	99
14) Pentachlorophenol	16.760	266	2135	0.776	ng/ul	97
15) Phenanthrene	17.149	178	12663	0.370	ng/ul	98
16) Anthracene	17.246	178	12367	0.385	ng/ul	99
19) Fluoranthene	19.168	202	14324	0.334	ng/ul	98
20) Pyrene	19.531	202	14784	0.337	ng/ul	98
21) Benzo(a)anthracene	21.289	228	12261	0.368	ng/ul	99
22) Chrysene	21.339	228	14899	0.398	ng/ul	99
24) Benzo(b)fluoranthene	22.873	252	14169	0.359	ng/ul	94
25) Benzo(k)fluoranthene	22.920	252	16576	0.388	ng/ul	94
26) Benzo(a)pyrene	23.455	252	13913	0.388	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	25.820	276	18392	0.446	ng/ul#	47
28) Dibenzo(a,h)anthracene	25.846	278	14486	0.456	ng/ul	99
29) Benzo(g,h,i)perylene	26.511	276	15930	0.440	ng/ul	99

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

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