

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032224\
 Data File : BM044716.D
 Acq On : 22 Mar 2024 01:57
 Operator : MA/JU
 Sample : PB159694BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS694

Quant Time: Mar 22 02:42:06 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	7.698	152	3499	0.400	ng/ul	0.00
4) Naphthalene-d8	10.479	136	10699	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.339	164	5120	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.102	188	10280	0.400	ng/ul	0.00
17) Chrysene-d12	21.298	240	8312	0.400	ng/ul	0.00
23) Perylene-d12	23.540	264	9354	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.260	96	3573	0.779	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.079	152	5598	0.364	ng/ul	0.00
18) Fluoranthene-d10	19.131	212	9664	0.331	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.289	88	10107	2.156	ng/ul#	64
5) Naphthalene	10.528	128	11396	0.379	ng/ul	98
7) 2-Methylnaphthalene	12.156	142	6625	0.348	ng/ul	99
8) 1-Methylnaphthalene	12.371	142	6815	0.354	ng/ul	100
10) Acenaphthylene	14.061	152	10296	0.396	ng/ul	97
11) Acenaphthene	14.404	153	7291	0.388	ng/ul	99
12) Fluorene	15.408	166	7962	0.371	ng/ul	99
14) Pentachlorophenol	16.756	266	2093	0.844	ng/ul	98
15) Phenanthrene	17.144	178	11737	0.380	ng/ul	98
16) Anthracene	17.241	178	11776	0.406	ng/ul	98
19) Fluoranthene	19.164	202	13455	0.331	ng/ul	98
20) Pyrene	19.526	202	13994	0.336	ng/ul	97
21) Benzo(a)anthracene	21.284	228	11390	0.360	ng/ul	99
22) Chrysene	21.333	228	14677	0.413	ng/ul	99
24) Benzo(b)fluoranthene	22.865	252	13472	0.377	ng/ul	93
25) Benzo(k)fluoranthene	22.908	252	15780	0.408	ng/ul	94
26) Benzo(a)pyrene	23.443	252	13154	0.406	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	25.803	276	17155	0.460	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.836	278	13625	0.474	ng/ul	99
29) Benzo(g,h,i)perylene	26.490	276	15137	0.462	ng/ul	99

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

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