

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110723\
 Data File : BM042630.D
 Acq On : 08 Nov 2023 01:57
 Operator : MA/JU
 Sample : PB156701BS
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS701

Quant Time: Nov 08 02:50:17 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 07 02:23:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	1293	0.400	ng/ul	0.00
4) Naphthalene-d8	10.677	136	3193	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.510	164	1903	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.266	188	3918	0.400	ng/ul	0.00
17) Chrysene-d12	21.447	240	2863	0.400	ng/ul	# 0.00
23) Perylene-d12	23.829	264	3321	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.206	96	1274	0.746	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.266	152	1438	0.323	ng/ul	0.00
18) Fluoranthene-d10	19.293	212	3428	0.346	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.240	88	3843	1.892	ng/ul#	58
5) Naphthalene	10.726	128	3448	0.334	ng/ul	99
7) 2-Methylnaphthalene	12.343	142	2176	0.329	ng/ul	95
8) 1-Methylnaphthalene	12.557	142	2154	0.321	ng/ul	99
10) Acenaphthylene	14.242	152	4191	0.322	ng/ul	98
11) Acenaphthene	14.575	153	2923	0.331	ng/ul	99
12) Fluorene	15.565	166	3329	0.328	ng/ul	100
14) Pentachlorophenol	16.908	266	1432	0.717	ng/ul	96
15) Phenanthrene	17.309	178	5069	0.334	ng/ul	99
16) Anthracene	17.406	178	4494	0.327	ng/ul	99
19) Fluoranthene	19.321	202	6014	0.336	ng/ul	100
20) Pyrene	19.689	202	6581	0.335	ng/ul	99
21) Benzo(a)anthracene	21.429	228	4195	0.329	ng/ul	100
22) Chrysene	21.485	228	4592	0.349	ng/ul	99
24) Benzo(b)fluoranthene	23.089	252	4847	0.361	ng/ul	95
25) Benzo(k)fluoranthene	23.142	252	5130	0.365	ng/ul	95
26) Benzo(a)pyrene	23.724	252	5205	0.375	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.316	276	6126	0.339	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.336	278	4514	0.350	ng/ul	94
29) Benzo(g,h,i)perylene	27.090	276	5552	0.331	ng/ul	98

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

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