

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091123\
 Data File : BM041791.D
 Acq On : 11 Sep 2023 22:39
 Operator : MA/JU
 Sample : PB155221BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS221

Quant Time: Sep 12 04:28:29 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM090523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 09 23:26:58 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.979	152	4484	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.796	136	10044	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.624	164	3937	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.375	188	7411	0.400	ng/ul	0.00
17) Chrysene-d12	21.545	240	2913	0.400	ng/ul	0.00
23) Perylene-d12	23.965	264	3101	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.389	96	4450	0.649	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.374	152	4793	0.371	ng/ul	-0.01
18) Fluoranthene-d10	19.394	212	5335	0.378	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.427	88	10734	1.508	ng/ul	91
5) Naphthalene	10.845	128	10991	0.333	ng/ul	100
7) 2-Methylnaphthalene	12.451	142	6175	0.330	ng/ul	100
8) 1-Methylnaphthalene	12.671	142	6539	0.348	ng/ul	99
10) Acenaphthylene	14.352	152	7988	0.328	ng/ul	99
11) Acenaphthene	14.685	153	6463	0.342	ng/ul	99
12) Fluorene	15.670	166	7100	0.335	ng/ul	99
14) Pentachlorophenol	16.999	266	2120	0.715	ng/ul	99
15) Phenanthrene	17.417	178	10976	0.351	ng/ul	99
16) Anthracene	17.506	178	9280	0.346	ng/ul	99
19) Fluoranthene	19.427	202	11520	0.335	ng/ul	99
20) Pyrene	19.789	202	11682	0.342	ng/ul	95
21) Benzo(a)anthracene	21.527	228	7944	0.334	ng/ul	100
22) Chrysene	21.583	228	8710	0.340	ng/ul	99
24) Benzo(b)fluoranthene	23.217	252	9208	0.332	ng/ul	95
25) Benzo(k)fluoranthene	23.266	252	8550	0.315	ng/ul	95
26) Benzo(a)pyrene	23.851	252	7434	0.330	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.471	276	10754	0.348	ng/ul#	92
28) Dibenzo(a,h)anthracene	26.488	278	7680	0.354	ng/ul	97
29) Benzo(g,h,i)perylene	27.250	276	9508	0.347	ng/ul	97

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

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