

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110921\
 Data File : BM032978.D
 Acq On : 11 Nov 2021 07:08
 Operator : CG/JU
 Sample : PB140354BS
 Misc :
 ALS Vial : 74 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS354

Quant Time: Nov 11 09:37:43 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM110921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 09 13:31:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.462	152	1930	0.400	ng/ul	0.00
4) Naphthalene-d8	10.235	136	7260	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.117	164	4159	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.858	188	8521	0.400	ng/ul	0.00
17) Chrysene-d12	21.049	240	6662	0.400	ng/ul	0.00
23) Perylene-d12	23.158	264	5154	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.860	96	1774	0.720	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.831	152	3525	0.342	ng/ul	0.00
18) Fluoranthene-d10	18.888	212	7650	0.379	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.895	88	5024	1.985	ng/ul#	89
5) Naphthalene	10.280	128	7659	0.363	ng/ul	99
7) 2-Methylnaphthalene	11.908	142	4798	0.329	ng/ul	100
8) 1-Methylnaphthalene	12.128	142	4594	0.321	ng/ul	100
10) Acenaphthylene	13.837	152	5920	0.327	ng/ul	99
11) Acenaphthene	14.178	153	5501	0.348	ng/ul	100
12) Fluorene	15.171	166	6141	0.330	ng/ul	99
14) Pentachlorophenol	16.532	266	1196	0.539	ng/ul	97
15) Phenanthrene	16.900	178	9666	0.341	ng/ul	99
16) Anthracene	16.990	178	7848	0.315	ng/ul	99
19) Fluoranthene	18.918	202	11154	0.357	ng/ul	98
20) Pyrene	19.284	202	11515	0.354	ng/ul	99
21) Benzo(a)anthracene	21.034	228	8103	0.329	ng/ul	99
22) Chrysene	21.085	228	9930	0.358	ng/ul	100
24) Benzo(b)fluoranthene	22.533	252	8643	0.366	ng/ul	98
25) Benzo(k)fluoranthene	22.574	252	9604	0.379	ng/ul	97
26) Benzo(a)pyrene	23.063	252	6996	0.359	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.215	276	8472	0.378	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.217	278	6750	0.380	ng/ul	98
29) Benzo(g,h,i)perylene	25.845	276	7672	0.394	ng/ul	98

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

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