

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111921\
 Data File : BM033201.D
 Acq On : 20 Nov 2021 06:49
 Operator : CG/JU
 Sample : PB140882BS
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS882

Quant Time: Nov 22 00:27:50 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 19 15:41:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.960	152	2653	0.400	ng/ul	0.00
4) Naphthalene-d8	10.756	136	7299	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.576	164	4425	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.313	188	9765	0.400	ng/ul	0.00
17) Chrysene-d12	21.466	240	7483	0.400	ng/ul	0.00
23) Perylene-d12	23.803	264	6038	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.417	96	1602	0.642	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.339	152	3691	0.365	ng/ul	0.00
18) Fluoranthene-d10	19.327	212	9100	0.420	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.448	88	4755	1.695	ng/ul	93
5) Naphthalene	10.805	128	7942	0.351	ng/ul	99
7) 2-Methylnaphthalene	12.411	142	5233	0.358	ng/ul	99
8) 1-Methylnaphthalene	12.627	142	5031	0.347	ng/ul	98
10) Acenaphthylene	14.299	152	7425	0.359	ng/ul#	99
11) Acenaphthene	14.641	153	6045	0.356	ng/ul	98
12) Fluorene	15.623	166	6963	0.364	ng/ul	98
14) Pentachlorophenol	16.959	266	1737	0.533	ng/ul	100
15) Phenanthrene	17.355	178	11259	0.318	ng/ul	99
16) Anthracene	17.445	178	9955	0.348	ng/ul	99
19) Fluoranthene	19.357	202	12767	0.354	ng/ul	98
20) Pyrene	19.720	202	13009	0.364	ng/ul	96
21) Benzo(a)anthracene	21.449	228	10041	0.376	ng/ul	100
22) Chrysene	21.502	228	11089	0.377	ng/ul	99
24) Benzo(b)fluoranthene	23.096	252	10521	0.385	ng/ul	99
25) Benzo(k)fluoranthene	23.142	252	10150	0.377	ng/ul	98
26) Benzo(a)pyrene	23.701	252	8489	0.375	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.197	276	10824	0.379	ng/ul#	91
28) Dibenzo(a,h)anthracene	26.216	278	8574	0.376	ng/ul	96
29) Benzo(g,h,i)perylene	26.931	276	9584	0.374	ng/ul	95

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

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