

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012622\
 Data File : BM033878.D
 Acq On : 26 Jan 2022 19:46
 Operator : CG/JU
 Sample : PB142272BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS272

Quant Time: Jan 27 01:42:52 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM012422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 24 15:20:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.939	152	1665	0.400	ng/ul	0.00
4) Naphthalene-d8	10.751	136	5683	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.579	164	2660	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	5379	0.400	ng/ul	0.00
17) Chrysene-d12	21.477	240	4646	0.400	ng/ul #	0.00
23) Perylene-d12	23.875	264	4623	0.400	ng/ul #-	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.296	96	1321	0.759	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.344	152	2995	0.382	ng/ul	0.00
18) Fluoranthene-d10	19.334	212	6048	0.423	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.330	88	3461	1.900	ng/ul#	78
5) Naphthalene	10.801	128	6002	0.384	ng/ul#	90
7) 2-Methylnaphthalene	12.415	142	3575	0.352	ng/ul	97
8) 1-Methylnaphthalene	12.631	142	3574	0.359	ng/ul	98
10) Acenaphthylene	14.303	152	5022	0.422	ng/ul#	93
11) Acenaphthene	14.640	153	3686	0.387	ng/ul	97
12) Fluorene	15.626	166	4196	0.386	ng/ul	98
14) Pentachlorophenol	16.969	266	1652	0.668	ng/ul	98
15) Phenanthrene	17.358	178	6845	0.399	ng/ul	97
16) Anthracene	17.451	178	6141	0.383	ng/ul	94
19) Fluoranthene	19.364	202	8658	0.409	ng/ul	99
20) Pyrene	19.723	202	8963	0.413	ng/ul#	94
21) Benzo(a)anthracene	21.462	228	7608	0.401	ng/ul	98
22) Chrysene	21.513	228	8884	0.441	ng/ul	97
24) Benzo(b)fluoranthene	23.141	252	8685	0.418	ng/ul	93
25) Benzo(k)fluoranthene	23.187	252	9183	0.441	ng/ul#	93
26) Benzo(a)pyrene	23.766	252	8447	0.469	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	26.383	276	10132	0.423	ng/ul#	85
28) Dibenzo(a,h)anthracene	26.404	278	8106	0.429	ng/ul	95
29) Benzo(g,h,i)perylene	27.138	276	9022	0.444	ng/ul	94

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

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