

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072621\
 Data File : BM031167.D
 Acq On : 26 Jul 2021 17:41
 Operator : CG/JU
 Sample : PB137838BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS838

Quant Time: Jul 26 18:11:21 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 26 17:09:27 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.632	152	1282	0.400	ng/ul	0.00
4) Naphthalene-d8	10.397	136	4983	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.262	164	3308	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.006	188	7277	0.400	ng/ul	0.00
17) Chrysene-d12	21.193	240	5579	0.400	ng/ul	0.00
23) Perylene-d12	23.363	264	4185	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	913	0.678	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.993	152	3041	0.348	ng/ul	0.00
18) Fluoranthene-d10	19.035	212	7276	0.428	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.248	88	2481	1.885	ng/ul#	83
5) Naphthalene	10.446	128	5150	0.346	ng/ul	99
7) 2-Methylnaphthalene	12.065	142	3664	0.342	ng/ul	100
8) 1-Methylnaphthalene	12.286	142	3391	0.324	ng/ul	100
10) Acenaphthylene	13.981	152	4690	0.347	ng/ul	98
11) Acenaphthene	14.322	153	4053	0.339	ng/ul	99
12) Fluorene	15.312	166	4757	0.338	ng/ul	99
14) Pentachlorophenol	16.659	266	1429	0.567	ng/ul	97
15) Phenanthrene	17.048	178	7988	0.346	ng/ul	99
16) Anthracene	17.138	178	7386	0.341	ng/ul	99
19) Fluoranthene	19.066	202	9158	0.408	ng/ul	98
20) Pyrene	19.428	202	9064	0.395	ng/ul	99
21) Benzo(a)anthracene	21.176	228	8081	0.363	ng/ul	99
22) Chrysene	21.227	228	8516	0.353	ng/ul	99
24) Benzo(b)fluoranthene	22.718	252	7476	0.380	ng/ul	99
25) Benzo(k)fluoranthene	22.757	252	7676	0.384	ng/ul	98
26) Benzo(a)pyrene	23.270	252	6184	0.364	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	25.516	276	7330	0.329	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.533	278	6017	0.337	ng/ul	99
29) Benzo(g,h,i)perylene	26.178	276	6121	0.324	ng/ul	97

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

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