

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080921\
 Data File : BM031604.D
 Acq On : 09 Aug 2021 15:32
 Operator : CG/JU
 Sample : M3244-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JCE01

Quant Time: Aug 09 16:27:35 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 09 10:16:43 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.580	152	1180	0.400	ng/ul	0.00
4) Naphthalene-d8	10.337	136	4475	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.212	164	2665	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.960	188	5150	0.400	ng/ul #	0.00
17) Chrysene-d12	21.158	240	4006	0.400	ng/ul #	0.00
23) Perylene-d12	23.319	264	3747	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.189	96	1838	1.402	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.934	152	1737	0.230	ng/ul	0.00
18) Fluoranthene-d10	18.997	212	3292	0.246	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.387	128	758	0.057	ng/ul#	77
7) 2-Methylnaphthalene	12.010	142	310	0.034	ng/ul	94
8) 1-Methylnaphthalene	12.235	142	192	0.021	ng/ul	95
14) Pentachlorophenol	16.620	266	44	0.027	ng/ul	89
15) Phenanthrene	17.002	178	1048	0.064	ng/ul#	78
19) Fluoranthene	19.027	202	1443	0.084	ng/ul#	32
20) Pyrene	19.389	202	1135	0.065	ng/ul#	26
21) Benzo(a)anthracene	21.142	228	558	0.034	ng/ul#	45
22) Chrysene	21.195	228	1714	0.102	ng/ul#	84
24) Benzo(b)fluoranthene	22.677	252	1299	0.077	ng/ul#	1
25) Benzo(k)fluoranthene	22.723	252	683	0.039	ng/ul#	1
26) Benzo(a)pyrene	23.222	252	484	0.033	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	25.456	276	541	0.028	ng/ul#	57

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

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