

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080521\
 Data File : BM031495.D
 Acq On : 05 Aug 2021 23:13
 Operator : CG/JU
 Sample : M3162-06
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGET6

Manual Integrations
 APPROVED

mohammad
 8/6/2021 2:49:16 PM

Quant Time: Aug 06 01:40:35 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 : Thu Aug 05 11:03:39 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.590	152	1294	0.400	ng/ul	0.00
4) Naphthalene-d8	10.347	136	4984	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.220	164	3168	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.968	188	5920	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.161	240	4601	0.400	ng/ul	0.00
23) Perylene-d12	23.316	264	4842	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.185	96	6661	4.635	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.944	152	2807	0.334	ng/ul	0.00
18) Fluoranthene-d10	19.001	212	5644	0.368	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.009	178	1615	0.086	ng/ul#	93
19) Fluoranthene	19.031	202	4214	0.214	ng/ul#	93
20) Pyrene	19.393	202	3585m	0.180	ng/ul	
21) Benzo(a)anthracene	21.147	228	1522	0.082	ng/ul#	92
22) Chrysene	21.197	228	1967	0.102	ng/ul	96
24) Benzo(b)fluoranthene	22.679	252	2725m	0.125	ng/ul	
25) Benzo(k)fluoranthene	22.716	252	1036m	0.046	ng/ul	
26) Benzo(a)pyrene	23.222	252	1688	0.088	ng/ul#	45
27) Indeno(1,2,3-cd)pyrene	25.458	276	1194	0.048	ng/ul#	100
29) Benzo(g,h,i)perylene	26.110	276	1218	0.058	ng/ul#	76

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

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