

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080521\
 Data File : BM031503.D
 Acq On : 06 Aug 2021 04:44
 Operator : CG/JU
 Sample : M3162-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGET1

Manual Integrations
 APPROVED

mohammad
 8/6/2021 2:49:35 PM

Quant Time: Aug 06 06:45:59 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 : Fri Aug 06 01:41:10 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	1183	0.400	ng/ul	0.00
4) Naphthalene-d8	10.347	136	4597	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.220	164	3064	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.968	188	5720	0.400	ng/ul	0.00
17) Chrysene-d12	21.159	240	4344	0.400	ng/ul	# 0.00
23) Perylene-d12	23.314	264	4575	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.186	96	9491	7.224	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.944	152	3944	0.509	ng/ul	0.00
18) Fluoranthene-d10	19.001	212	7983	0.551	ng/ul	0.00
Target Compounds						
					Qvalue	
10) Acenaphthylene	13.941	152	308	0.024	ng/ul#	53
15) Phenanthrene	17.009	178	1854	0.102	ng/ul#	93
16) Anthracene	17.100	178	463	0.027	ng/ul#	75
19) Fluoranthene	19.028	202	5050	0.272	ng/ul#	93
20) Pyrene	19.394	202	4673m	0.248	ng/ul	
21) Benzo(a)anthracene	21.144	228	2508	0.143	ng/ul	94
22) Chrysene	21.195	228	2918m	0.161	ng/ul	
24) Benzo(b)fluoranthene	22.675	252	4222m	0.205	ng/ul	
25) Benzo(k)fluoranthene	22.713	252	1359m	0.064	ng/ul	
26) Benzo(a)pyrene	23.222	252	2711	0.150	ng/ul#	68
27) Indeno(1,2,3-cd)pyrene	25.454	276	1867	0.080	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.458	278	504	0.027	ng/ul#	1
29) Benzo(g,h,i)perylene	26.105	276	1906	0.096	ng/ul#	83

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

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