

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062521\
 Data File : BM030615.D
 Acq On : 25 Jun 2021 23:21
 Operator : CG/JU
 Sample : M2682-05DL 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleID :
 BGD76DL

Manual Integrations
 APPROVED

mohammad
 6/28/2021 1:21:50 PM

Quant Time: Jun 26 01:32:12 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 25 10:09:59 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	4114	0.400	ng/ul	0.00
4) Naphthalene-d8	10.586	136	16782	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.425	164	9822	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.159	188	18008	0.400	ng/ul	0.00
17) Chrysene-d12	21.326	240	11494	0.400	ng/ul	0.00
23) Perylene-d12	23.581	264	11513	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.282	96	1583	0.360	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.178	152	685	0.026	ng/ul	0.00
18) Fluoranthene-d10	19.180	212	1293	0.037	ng/ul	0.00
Target Compounds						
					Qvalue	
14) Pentachlorophenol	16.832	266	119	0.023	ng/ul	95
15) Phenanthrene	17.200	178	6172	0.096	ng/ul	95
16) Anthracene	17.291	178	1375	0.024	ng/ul#	83
19) Fluoranthene	19.211	202	16154	0.306	ng/ul	99
20) Pyrene	19.573	202	10938	0.202	ng/ul	100
21) Benzo(a)anthracene	21.311	228	5975	0.131	ng/ul#	88
22) Chrysene	21.362	228	5725	0.113	ng/ul#	91
24) Benzo(b)fluoranthene	22.902	252	7049m	0.141	ng/ul	
25) Benzo(k)fluoranthene	22.946	252	2205m	0.042	ng/ul	
26) Benzo(a)pyrene	23.481	252	2244	0.050	ng/ul#	23
27) Indeno(1,2,3-cd)pyrene	25.860	276	1916	0.034	ng/ul#	100

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

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