

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080221\
 Data File : BM031456.D
 Acq On : 04 Aug 2021 20:25
 Operator : CG/JU
 Sample : M3141-18DL 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGER9DL

Manual Integrations
 APPROVED

mohammad
 8/5/2021 5:04:45 PM

Quant Time: Aug 05 01:01:52 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 : Wed Aug 04 12:17:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.594	152	1612	0.400	ng/ul	0.00
4) Naphthalene-d8	10.356	136	6090	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.224	164	3647	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.971	188	7088	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.164	240	5773	0.400	ng/ul	0.00
23) Perylene-d12	23.321	264	5491	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.192	96	3915	2.187	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.948	152	1375	0.134	ng/ul	0.00
18) Fluoranthene-d10	19.005	212	2554	0.133	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.401	128	689	0.038	ng/ul#	79
14) Pentachlorophenol	16.631	266	201	0.089	ng/ul	99
15) Phenanthrene	17.013	178	1770	0.079	ng/ul	94
16) Anthracene	17.103	178	1041	0.049	ng/ul#	85
19) Fluoranthene	19.035	202	14293	0.580	ng/ul	99
20) Pyrene	19.397	202	14662	0.586	ng/ul	98
21) Benzo(a)anthracene	21.149	228	11184	0.480	ng/ul	94
22) Chrysene	21.200	228	10040	0.416	ng/ul	98
24) Benzo(b)fluoranthene	22.682	252	11956m	0.483	ng/ul	
25) Benzo(k)fluoranthene	22.720	252	4585m	0.179	ng/ul	
26) Benzo(a)pyrene	23.227	252	10174	0.470	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.458	276	5244	0.187	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.463	278	1460	0.065	ng/ul#	61
29) Benzo(g,h,i)perylene	26.115	276	4684	0.196	ng/ul	97

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

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