

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070821\
 Data File : BN015479.D
 Acq On : 10 Jul 2021 12:12
 Operator : CG/JU
 Sample : M2808-10DL 5X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGDR8DL

Manual Integrations
 APPROVED

mohammad
 7/12/2021 9:34:25 AM

Quant Time: Jul 11 06:14:56 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 08 11:17:31 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.711	152	3569	0.400	ng/ul	0.00
4) Naphthalene-d8	10.480	136	12188	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.335	164	6252	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.082	188	12310	0.400	ng/ul	0.00
17) Chrysene-d12	21.281	240	9620	0.400	ng/ul	0.00
23) Perylene-d12	23.520	264	8669	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	1021	0.254	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.069	152	438	0.023	ng/ul	0.00
18) Fluoranthene-d10	19.118	212	880	0.028	ng/ul	0.00
Target Compounds						
					Qvalue	
11) Acenaphthene	14.395	153	574	0.023	ng/ul	96
15) Phenanthrene	17.124	178	6959	0.156	ng/ul	98
16) Anthracene	17.213	178	1511	0.038	ng/ul#	90
19) Fluoranthene	19.146	202	15909	0.358	ng/ul	98
20) Pyrene	19.513	202	12314	0.276	ng/ul	99
21) Benzo(a)anthracene	21.267	228	5696	0.160	ng/ul	96
22) Chrysene	21.319	228	5920	0.139	ng/ul	98
24) Benzo(b)fluoranthene	22.851	252	7401m	0.183	ng/ul	
25) Benzo(k)fluoranthene	22.889	252	2944m	0.064	ng/ul	
26) Benzo(a)pyrene	23.421	252	4338	0.128	ng/ul#	74
27) Indeno(1,2,3-cd)pyrene	25.760	276	3199	0.078	ng/ul#	89
28) Dibenzo(a,h)anthracene	25.770	278	768	0.024	ng/ul#	54

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

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