

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071221\
 Data File : BN015552.D
 Acq On : 14 Jul 2021 05:55
 Operator : CG/JU
 Sample : M2869-11DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Jul 14 07:17:49 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 : Wed Jul 14 07:10:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	3534	0.40	ng/ul	0.00
4) Naphthalene-d8	10.595	136	12518	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.441	164	6561	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.191	188	13008	0.40	ng/ul	0.00
17) Chrysene-d12	21.384	240	11008	0.40	ng/ul	0.00
23) Perylene-d12	23.736	264	9902	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	1019	0.26	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.190	152	281	0.01	ng/ul	0.00
18) Fluoranthene-d10	19.225	212	602	0.02	ng/ul	0.00
Target Compounds						
						Qvalue
5) Naphthalene	10.644	128	2049	0.05	ng/ul#	93
10) Acenaphthylene	14.159	152	5408	0.16	ng/ul	99
11) Acenaphthene	14.506	153	1201	0.05	ng/ul	98
12) Fluorene	15.491	166	2071	0.07	ng/ul#	95
15) Phenanthrene	17.234	178	59182	1.25	ng/ul	100
16) Anthracene	17.322	178	32325	0.78	ng/ul	100
19) Fluoranthene	19.253	202	176197	3.46	ng/ul	98
20) Pyrene	19.615	202	119467	2.34	ng/ul	98
21) Benzo(a)anthracene	21.369	228	85671	2.10	ng/ul	95
22) Chrysene	21.419	228	66597	1.36	ng/ul	98
24) Benzo(b)fluoranthene	23.026	252	79256m	1.71	ng/ul	
25) Benzo(k)fluoranthene	23.067	252	29371m	0.56	ng/ul	
26) Benzo(a)pyrene	23.631	252	36550	0.94	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.156	276	24610	0.52	ng/ul#	92
28) Dibenzo(a,h)anthracene	26.172	278	8100	0.22	ng/ul	92
29) Benzo(g,h,i)perylene	26.897	276	1383m	0.03	ng/ul	

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

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