

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071221\
 Data File : BN015549.D
 Acq On : 14 Jul 2021 04:07
 Operator : CG/JU
 Sample : M2869-09
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGDx7

Manual Integrations
 APPROVED

mohammad
 7/14/2021 12:15:14 PM

Quant Time: Jul 14 05:42:52 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 05:24:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.808	152	3057	0.400	ng/ul	0.00
4) Naphthalene-d8	10.595	136	10837	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.441	164	5971	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.191	188	11701m	0.400	ng/ul	0.00
17) Chrysene-d12	21.384	240	9503	0.400	ng/ul	0.00
23) Perylene-d12	23.739	264	9073	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.310	96	13789	4.002	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.184	152	4537	0.267	ng/ul	0.00
18) Fluoranthene-d10	19.225	212	9219	0.295	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.644	128	791	0.023	ng/ul#	77
10) Acenaphthylene	14.159	152	1061	0.034	ng/ul#	67
11) Acenaphthene	14.506	153	37054	1.566	ng/ul	98
12) Fluorene	15.492	166	2189	0.084	ng/ul#	97
15) Phenanthrene	17.234	178	42147	0.992	ng/ul	100
16) Anthracene	17.322	178	7644	0.204	ng/ul	99
19) Fluoranthene	19.253	202	19704	0.449	ng/ul	98
20) Pyrene	19.615	202	14969	0.340	ng/ul	96
21) Benzo(a)anthracene	21.369	228	6324	0.180	ng/ul	93
22) Chrysene	21.422	228	4789	0.114	ng/ul	97
24) Benzo(b)fluoranthene	23.029	252	5524m	0.130	ng/ul	
25) Benzo(k)fluoranthene	23.070	252	2283m	0.048	ng/ul	
26) Benzo(a)pyrene	23.637	252	4236	0.119	ng/ul#	80
27) Indeno(1,2,3-cd)pyrene	26.159	276	2278	0.053	ng/ul#	98
29) Benzo(g,h,i)perylene	26.900	276	1911	0.048	ng/ul#	85

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

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