

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072920\
 Data File : BN011290.D
 Acq On : 29 Jul 2020 13:21
 Operator : CG/JU
 Sample : L3388-02DL 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD0DL

Manual Integrations
 APPROVED

mohammad
 7/30/2020 3:35:26 PM

Quant Time: Jul 29 14:01:07 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN072020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 29 11:20:12 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	1452	0.40	ng/ul	0.00
2) Naphthalene-d8	10.22	136	5267	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	3203	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	7370	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	6961	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	6321	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	400	0.03	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	1125	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.27	128	800	0.051	ng/ul#	84
5) 2-Methylnaphthalene	11.91	142	243	0.024	ng/ul	83
7) Acenaphthylene	13.82	152	487	0.031	ng/ul#	72
8) Acenaphthene	14.17	153	611	0.050	ng/ul	96
9) Fluorene	15.17	166	582	0.042	ng/ul#	96
12) Phenanthrene	16.90	178	12534	0.516	ng/ul	99
13) Anthracene	17.00	178	2593	0.130	ng/ul	96
15) Fluoranthene	18.93	202	27539	1.015	ng/ul	99
17) Pyrene	19.30	202	24211	0.651	ng/ul	100
18) Benzo(a)anthracene	21.07	228	15193	0.617	ng/ul	98
19) Chrysene	21.12	228	15203	0.464	ng/ul	98
21) Benzo(b)fluoranthene	22.59	252	19108m	0.700	ng/ul	
22) Benzo(k)fluoranthene	22.62	252	7513m	0.240	ng/ul	
23) Benzo(a)pyrene	23.12	252	10781	0.470	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.29	276	7048	0.260	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.29	278	2042	0.100	ng/ul#	78
26) Benzo(a,h,i)perylene	25.91	276	4491	0.167	ng/ul	95

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

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