

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072920\
 Data File : BN011289.D
 Acq On : 29 Jul 2020 10:21
 Operator : CG/JU
 Sample : L3388-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD0

Manual Integrations
 APPROVED

mohammad
 7/30/2020 3:35:23 PM

Quant Time: Jul 29 11:28:15 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	1423	0.40	ng/ul	0.00
2) Naphthalene-d8	10.22	136	4807	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	2890	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.87	188	7090m	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	6046	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	4846	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.82	152	1708	0.15	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	4684	0.15	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.27	128	2904	0.202	ng/ul	99
5) 2-Methylnaphthalene	11.90	142	731	0.078	ng/ul	99
7) Acenaphthylene	13.82	152	1216	0.087	ng/ul#	89
8) Acenaphthene	14.17	153	2238	0.203	ng/ul	99
9) Fluorene	15.17	166	2001	0.160	ng/ul#	99
12) Phenanthrene	16.90	178	52963	2.266	ng/ul	98
13) Anthracene	17.00	178	10352	0.538	ng/ul	98
15) Fluoranthene	18.93	202	114256	4.376	ng/ul	99
17) Pyrene	19.30	202	93794	2.906	ng/ul	100
18) Benzo(a)anthracene	21.06	228	57139	2.673	ng/ul	98
19) Chrysene	21.11	228	54075	1.902	ng/ul	98
21) Benzo(b)fluoranthene	22.58	252	66688m	3.187	ng/ul	
22) Benzo(k)fluoranthene	22.62	252	25761m	1.073	ng/ul	
23) Benzo(a)pyrene	23.12	252	37638	2.138	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	25.28	276	26480	1.277	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.29	278	7597	0.486	ng/ul	94
26) Benzo(a,h,i)perylene	25.91	276	17940	0.870	ng/ul	98

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

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