

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121119\
 Data File : BN008945.D
 Acq On : 12 Dec 2019 14:53
 Operator : JU
 Sample : K6089-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BP9

Manual Integrations
 APPROVED

mohammad
 12/13/2019 9:07:33 AM

Quant Time: Dec 12 17:25:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 12 00:49:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2567	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	10845	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	6962	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	15784	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	13826	0.40	ng/ul	0.00
20) Perylene-d12	23.27	264	14827	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	1616	0.09	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	4556	0.09	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.36	128	2379	0.073	ng/ul#	92
5) 2-Methylnaphthalene	11.98	142	2129	0.093	ng/ul	99
7) Acenaphthylene	13.90	152	4654	0.121	ng/ul	96
9) Fluorene	15.24	166	879	0.026	ng/ul#	82
12) Phenanthrene	16.97	178	51518	0.968	ng/ul	99
13) Anthracene	17.06	178	3270	0.065	ng/ul#	93
15) Fluoranthene	18.99	202	82676	1.264	ng/ul	99
17) Pyrene	19.35	202	75537	1.362	ng/ul	99
18) Benzo(a)anthracene	21.11	228	35740	0.643	ng/ul	96
19) Chrysene	21.17	228	60190	1.102	ng/ul	98
21) Benzo(b)fluoranthene	22.64	252	68687m	1.123	ng/ul	
22) Benzo(k)fluoranthene	22.68	252	26537m	0.438	ng/ul	
23) Benzo(a)pyrene	23.18	252	31429	0.548	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.38	276	24310	0.362	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.39	278	6087	0.112	ng/ul#	83
26) Benzo(g,h,i)perylene	26.02	276	23538	0.419	ng/ul	99

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

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