

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121119\
 Data File : BN008929.D
 Acq On : 11 Dec 2019 22:21
 Operator : JU
 Sample : K6088-14
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BP2

Manual Integrations
 APPROVED

mohammad
 12/12/2019 11:32:33 AM

Quant Time: Dec 12 01:48:22 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 : Wed Dec 11 10:35:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2364	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	10116	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	6587	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	15346	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	13472	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	14256	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	2835	0.17	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	8891	0.19	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.97	178	2838	0.055	ng/ul#	92
15) Fluoranthene	19.00	202	9333	0.147	ng/ul	99
17) Pyrene	19.36	202	8430m	0.156	ng/ul	
18) Benzo(a)anthracene	21.12	228	4420	0.082	ng/ul	97
19) Chrysene	21.17	228	4705	0.088	ng/ul	96
21) Benzo(b)fluoranthene	22.64	252	5766m	0.098	ng/ul	
22) Benzo(k)fluoranthene	22.69	252	2761m	0.047	ng/ul	
23) Benzo(a)pyrene	23.19	252	4128	0.075	ng/ul#	70
24) Indeno(1,2,3-cd)pyrene	25.39	276	3032	0.047	ng/ul#	92
26) Benzo(g,h,i)perylene	26.02	276	2862	0.053	ng/ul#	89

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

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