

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082819\
 Data File : BN007455.D
 Acq On : 28 Aug 2019 05:06
 Operator : HP/JU
 Sample : K4337-10DL 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA0DL

Manual Integrations
 APPROVED

mohammad
 8/28/2019 7:49:03 AM

Quant Time: Aug 28 08:49:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 28 02:09:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	5600	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	23747	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	13425	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	28603	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	25076	0.40	ng/ul	0.00
20) Perylene-d12	23.12	264	28282	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	3056	0.08	ng/ul	0.00
14) Fluoranthene-d10	18.84	212	8657	0.09	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.20	128	3697	0.055	ng/ul#	92
7) Acenaphthylene	13.76	152	6126	0.082	ng/ul#	94
8) Acenaphthene	14.11	153	21448	0.400	ng/ul	98
9) Fluorene	15.11	166	10263	0.177	ng/ul#	94
11) Pentachlorophenol	16.46	266	20553	3.190	ng/ul	99
12) Phenanthrene	16.84	178	9017	0.101	ng/ul	96
13) Anthracene	16.93	178	30063m	0.357	ng/ul	
15) Fluoranthene	18.87	202	31420	0.263	ng/ul	99
17) Pyrene	19.24	202	26275	0.214	ng/ul	100
18) Benzo(a)anthracene	21.00	228	14961	0.133	ng/ul#	90
19) Chrysene	21.05	228	19367m	0.161	ng/ul	
21) Benzo(b)fluoranthene	22.50	252	34925m	0.283	ng/ul	
22) Benzo(k)fluoranthene	22.54	252	10630m	0.082	ng/ul	
23) Benzo(a)pyrene	23.03	252	16982	0.138	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.18	276	16547	0.120	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.19	278	3718	0.034	ng/ul#	58
26) Benzo(a,h,i)perylene	25.81	276	15178	0.129	ng/ul	97

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

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