

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090819\
 Data File : BN007568.D
 Acq On : 09 Sep 2019 01:49
 Operator : HP/JU
 Sample : K4762-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 STOCK-PILE

Manual Integrations
 APPROVED

mohammad
 9/16/2019 8:53:24 AM

Quant Time: Sep 09 13:30:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 29 12:18:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	7086	0.40	ng	-0.02
7) Naphthalene-d8	10.14	136	38319	0.40	ng	-0.03
13) Acenaphthene-d10	14.03	164	33946	0.40	ng	-0.02
19) Phenanthrene-d10	16.78	188	31413	0.40	ng	-0.02
27) Chrysene-d12	21.00	240	30396	0.40	ng	-0.01
34) Perylene-d12	23.11	264	35670	0.40	ng	-0.02

System Monitoring Compounds						
4) 2-Fluorophenol	5.02	112	8147	0.33	ng	-0.02
5) Phenol-d6	6.58	99	10808	0.34	ng	0.00
8) Nitrobenzene-d5	8.52	82	12938	0.39	ng	-0.01
11) 2-Methylnaphthalene-d10	11.75	152	23159	0.36	ng	-0.01
14) 2,4,6-Tribromophenol	15.54	330	2629	0.15	ng	-0.02
15) 2-Fluorobiphenyl	12.65	172	16820	0.12	ng	0.00
25) Fluoranthene-d10	18.83	212	31958	0.32	ng	-0.02
29) Terphenyl-d14	19.45	244	18838	0.24	ng	-0.02

Target Compounds				Qvalue		
9) Naphthalene	10.19	128	14548	0.133	ng	# 85
12) 2-Methylnaphthalene	11.82	142	9866	0.132	ng	# 24
16) Acenaphthylene	13.78	152	124525	0.620	ng	# 58
17) Acenaphthene	14.10	154	29109	0.248	ng	# 86
18) Fluorene	15.09	166	36876	0.248	ng	# 67
23) Phenanthrene	16.83	178	430804	4.559	ng	# 94
24) Anthracene	16.92	178	115005	1.210	ng	# 93
26) Fluoranthene	18.86	202	956104	7.881	ng	# 98
28) Pyrene	19.23	202	1122296	9.179	ng	# 97
30) Benzo(a)anthracene	20.99	228	545157	4.567	ng	# 87
31) Chrysene	21.04	228	524673	4.554	ng	# 90
32) Bis(2-ethylhexyl)phthalate	20.95	149	72238	0.916	ng	# 97
33) Indeno(1,2,3-cd)pyrene	25.18	276	335549	2.417	ng	# 99
35) Benzo(b)fluoranthene	22.49	252	701718	5.430	ng	# 90
36) Benzo(k)fluoranthene	22.53	252	223534m	1.750	ng	# 92
37) Benzo(a)pyrene	23.02	252	580426	4.719	ng	# 92
38) Dibenzo(a,h)anthracene	25.18	278	103978	0.847	ng	# 30
39) Benzo(a,h,i)perylene	25.81	276	425616	3.481	ng	# 90

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

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