

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121417\
 Data File : BM013430.D
 Acq On : 15 Dec 2017 08:23
 Operator : SJ/JU
 Sample : I6776-03 10X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F9A55

Manual Integrations
 APPROVED

Sohil
 12/15/2017 6:49:19 PM

Quant Time: Dec 25 01:36:33 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 15 04:56:09 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	259791	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	1116558	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.32	164	638396	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.07	188	1334302	20.00	ng/ul	0.00
75) Chrysene-d12	21.27	240	1128723	20.00	ng/ul	0.01
83) Perylene-d12	23.49	264	920773	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	1973	0.38	ng/uL	0.00
5) Phenol-d5	6.85	99	78010	3.47	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.02	67	45556	3.55	ng/ul	0.00
9) 2-Chlorophenol-d4	7.22	132	61628	3.51	ng/ul	0.00
13) 4-Methylphenol-d8	8.39	113	66218	3.45	ng/ul	0.00
19) Nitrobenzene-d5	8.83	128	32805	3.70	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	37423	3.95	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	68865	3.53	ng/ul	0.00
29) 4-Chloroaniline-d4	10.60	131	23251	1.30	ng/ul	0.00
43) Dimethylphthalate-d6	13.73	166	245538	4.32	ng/ul	0.00
46) Acenaphthylene-d8	14.01	160	296663	4.23	ng/ul	0.00
51) 4-Nitrophenol-d4	14.54	143	30865	3.13	ng/ul	0.00
57) Fluorene-d10	15.32	176	207965	4.26	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.46	200	23891	2.85	ng/ul	0.00
70) Anthracene-d10	17.16	188	305722	4.57	ng/ul	0.00
76) Pyrene-d10	19.48	212	319963	5.69	ng/ul	0.01
87) Benzo(a)pyrene-d12	23.35	264	201111	4.29	ng/ul	0.00

Target Compounds

					Ovalue
47) Acenaphthylene	14.04	152	573854	8.845	ng/ul 99
68) Pentachlorophenol	16.72	266	68921	6.889	ng/ul 94
69) Phenanthrene	17.11	178	661978	8.733	ng/ul 99
71) Anthracene	17.20	178	1276908	16.480	ng/ul 98
74) Fluoranthene	19.14	202	27754813m	298.773	ng/ul
77) Pyrene	19.51	202	31358470m	457.300	ng/ul
80) Benzo(a)anthracene	21.26	228	12193300	170.262	ng/ul 96
82) Chrysene	21.32	228	14542850m	218.886	ng/ul
85) Benzo(b)fluoranthene	22.84	252	8643897	139.259	ng/ul# 97
86) Benzo(k)fluoranthene	22.87	252	2908805	49.798	ng/ul# 98
88) Benzo(a)pyrene	23.40	252	2896149	51.941	ng/ul# 96
89) Indeno(1,2,3-cd)pyrene	25.72	276	1440219	26.032	ng/ul# 94
90) Dibenzo(a,h)anthracene	25.72	278	402854	8.552	ng/ul# 89
91) Benzo(g,h,i)perylene	26.40	276	1054633	24.167	ng/ul# 95

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

