

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP021721\
 Data File : BP004778.D
 Acq On : 17 Feb 2021 18:01
 Operator : CG/JU
 Sample : M1379-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBGY7

Manual Integrations
 APPROVED

mohammad
 2/18/2021 1:40:07 PM

Quant Time: Feb 18 07:20:42 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 17 23:17:49 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	49871	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	181035	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.41	164	103235	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.16	188	215323	20.00	ng/ul	0.00
78) Chrysene-d12	21.26	240	238148	20.00	ng/ul	0.00
86) Perylene-d12	23.56	264	248128	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	3880	3.01	ng/uL	0.00
5) Phenol-d5	6.93	99	64286	16.06	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.10	67	34875	16.78	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	53835	17.75	ng/ul	0.00
13) 4-Methylphenol-d8	8.47	113	47149	14.97	ng/ul	0.00
19) Nitrobenzene-d5	8.92	128	24940	19.95	ng/ul	0.00
22) 2-Nitrophenol-d4	9.64	143	27580	20.11	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.17	165	50889	18.43	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	22119	7.18	ng/ul	0.00
44) Dimethylphthalate-d6	13.82	166	157382	21.00	ng/ul	0.00
47) Acenaphthylene-d8	14.10	160	192806	21.15	ng/ul	0.00
52) 4-Nitrophenol-d4	14.62	143	18603	14.20	ng/ul	0.00
58) Fluorene-d10	15.40	176	137089	21.35	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.52	200	19753m	15.18	ng/ul	0.00
71) Anthracene-d10	17.26	188	215339	22.68	ng/ul	0.00
79) Pyrene-d10	19.51	212	249079	22.34	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.42	264	277158	22.08	ng/ul	0.00

Target Compounds

					Ovalue
4) Benzaldehyde	6.90	77	4571	2.180	ng/ul# 80
6) Phenol	6.96	94	7044m	1.688	ng/ul
14) Acetophenone	8.58	105	12376	2.389	ng/ul 99
45) Dimethylphthalate	13.87	163	12449	1.611	ng/ul# 99
77) Fluoranthene	19.18	202	22925	1.664	ng/ul# 92
80) Pyrene	19.53	202	21006	1.452	ng/ul 95
88) Benzo(b)fluoranthene	22.86	252	21836	1.398	ng/ul# 76

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

