

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP031220\
 Data File : BP002211.D
 Acq On : 13 Mar 2020 12:20
 Operator : CG/JU
 Sample : L1840-21
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBET6

Quant Time: Mar 13 14:15:45 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 13 03:14:27 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	246404	20.00	ng/ul	0.00
18) Naphthalene-d8	10.39	136	1029648	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.26	164	644397	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	1426320	20.00	ng/ul	0.00
78) Chrysene-d12	21.11	240	1431142	20.00	ng/ul	0.00
86) Perylene-d12	23.32	264	1702533	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.15	96	26627	4.65	ng/uL	0.00
5) Phenol-d5	6.81	99	64411	3.43	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.96	67	105193	10.57	ng/ul	0.00
9) 2-Chlorophenol-d4	7.16	132	137897	8.48	ng/ul	0.00
13) 4-Methylphenol-d8	8.33	113	94158	6.10	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	60135	7.60	ng/ul	0.00
22) 2-Nitrophenol-d4	9.49	143	61075	6.86	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	136656	7.97	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	12984	0.69	ng/ul	0.01
44) Dimethylphthalate-d6	13.67	166	468494	9.61	ng/ul	0.00
47) Acenaphthylene-d8	13.95	160	529107	9.07	ng/ul	0.00
52) 4-Nitrophenol-d4	14.49	143	12740	1.37	ng/ul	0.01
58) Fluorene-d10	15.26	176	428270	10.15	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	31449	4.00	ng/ul	0.00
71) Anthracene-d10	17.12	188	653717	10.20	ng/ul	0.00
79) Pyrene-d10	19.36	212	775045	10.35	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.17	264	839642	9.64	ng/ul	0.00

Target Compounds					Ovalue	
6) Phenol	6.83	94	65683	3.366	ng/ul	97
45) Dimethylphthalate	13.72	163	121155	2.407	ng/ul	99

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

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