

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP031220\
 Data File : BP002189.D
 Acq On : 12 Mar 2020 18:11
 Operator : CG/JU
 Sample : PB127498BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK98

Quant Time: Mar 13 02:54:07 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 12 13:49:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	248769	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	1030198	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.26	164	637640	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	1361001	20.00	ng/ul	0.00
78) Chrysene-d12	21.11	240	1341163	20.00	ng/ul	0.00
86) Perylene-d12	23.32	264	1564114	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.15	96	42341	7.32	ng/uL	0.00
5) Phenol-d5	6.81	99	736659	38.82	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.96	67	430041	42.80	ng/ul	0.00
9) 2-Chlorophenol-d4	7.16	132	643157	39.18	ng/ul	0.00
13) 4-Methylphenol-d8	8.33	113	595944	38.21	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	298749	37.75	ng/ul	0.00
22) 2-Nitrophenol-d4	9.49	143	329415	37.00	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	607948	35.45	ng/ul	0.00
29) 4-Chloroaniline-d4	10.53	131	861914	45.84	ng/ul	0.00
44) Dimethylphthalate-d6	13.68	166	1888836	39.14	ng/ul	0.00
47) Acenaphthylene-d8	13.95	160	2288317	39.65	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	273915	29.69	ng/ul	0.00
58) Fluorene-d10	15.26	176	1682428	40.29	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	210988	28.12	ng/ul	0.00
71) Anthracene-d10	17.12	188	2616980	42.78	ng/ul	0.00
79) Pyrene-d10	19.36	212	3021983	43.04	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.17	264	3367518	42.08	ng/ul	0.00

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

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