

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030521\
 Data File : BP004918.D
 Acq On : 05 Mar 2021 16:09
 Operator : CG/JU
 Sample : PB134929BL
 Misc : SOM-MDL-W-BL-05
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK29

Quant Time: Mar 05 16:41:31 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP030521MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 05 13:02:07 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	33555	20.00	ng/ul	0.00
18) Naphthalene-d8	10.53	136	129129	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.39	164	84034	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.15	188	185436	20.00	ng/ul	0.00
78) Chrysene-d12	21.24	240	196984	20.00	ng/ul	0.00
86) Perylene-d12	23.53	264	197808	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.23	96	5496	6.65	ng/uL	0.00
5) Phenol-d5	6.91	99	88524	34.51	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.08	67	48115	36.57	ng/ul	0.00
9) 2-Chlorophenol-d4	7.27	132	69275	33.87	ng/ul	0.00
13) 4-Methylphenol-d8	8.45	113	72323	34.89	ng/ul	0.00
19) Nitrobenzene-d5	8.89	128	33845	36.51	ng/ul	0.00
22) 2-Nitrophenol-d4	9.62	143	39600	37.39	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.15	165	68494	32.87	ng/ul	0.00
29) 4-Chloroaniline-d4	10.67	131	95638	42.04	ng/ul	0.00
44) Dimethylphthalate-d6	13.80	166	251253	41.56	ng/ul	0.00
47) Acenaphthylene-d8	14.08	160	278654	37.46	ng/ul	0.00
52) 4-Nitrophenol-d4	14.60	143	35994	33.91	ng/ul	0.00
58) Fluorene-d10	15.39	176	205919	38.22	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.51	200	39003	32.13	ng/ul	0.00
71) Anthracene-d10	17.25	188	346639	41.63	ng/ul	0.00
79) Pyrene-d10	19.49	212	402812	45.36	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.39	264	437062	43.48	ng/ul	0.00

Target Compounds Ovalue

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

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