

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030321\
 Data File : BP004895.D
 Acq On : 03 Mar 2021 16:01
 Operator : CG/JU
 Sample : PB134916BL
 Misc : SOM-MDL-S-BL02
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK16

Quant Time: Mar 03 16:33:08 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP030221MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 03 10:35:24 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	30017	20.00	ng/ul	0.00
18) Naphthalene-d8	10.53	136	118553	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.39	164	76653	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.15	188	167753	20.00	ng/ul	-0.01
78) Chrysene-d12	21.24	240	189120	20.00	ng/ul	-0.01
86) Perylene-d12	23.53	264	193347	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	5191	6.80	ng/uL	0.00
5) Phenol-d5	6.92	99	79058	32.34	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.08	67	44316	34.96	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.28	132	65696	34.84	ng/ul	0.00
13) 4-Methylphenol-d8	8.45	113	63834	32.77	ng/ul	-0.01
19) Nitrobenzene-d5	8.90	128	30549	35.38	ng/ul	0.00
22) 2-Nitrophenol-d4	9.62	143	35189	36.29	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	10.16	165	60388	32.93	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.68	131	83327	39.76	ng/ul	0.00
44) Dimethylphthalate-d6	13.80	166	211694	37.75	ng/ul	-0.01
47) Acenaphthylene-d8	14.09	160	240793	35.07	ng/ul	0.00
52) 4-Nitrophenol-d4	14.60	143	32696	32.59	ng/ul	-0.01
58) Fluorene-d10	15.39	176	176077	36.93	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.51	200	36375	33.25	ng/ul	-0.01
71) Anthracene-d10	17.25	188	294777	38.75	ng/ul	0.00
79) Pyrene-d10	19.49	212	349044	39.11	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.39	264	391499	39.51	ng/ul	-0.01

Target Compounds Ovalue

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

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