

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030321\
 Data File : BP004901.D
 Acq On : 03 Mar 2021 19:24
 Operator : CG/JU
 Sample : PB134917BL
 Misc : SOM-MDL-MES-BL03
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK17

Quant Time: Mar 04 03:06:44 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	28589	20.00	ng/ul	0.00
18) Naphthalene-d8	10.53	136	110793	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.39	164	73760	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.15	188	163955	20.00	ng/ul	-0.01
78) Chrysene-d12	21.24	240	183356	20.00	ng/ul	-0.01
86) Perylene-d12	23.53	264	187684	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.23	96	4350	5.99	ng/uL	0.00
5) Phenol-d5	6.92	99	73706	31.66	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.08	67	40957	33.92	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.28	132	60009	33.41	ng/ul	0.00
13) 4-Methylphenol-d8	8.45	113	59306	31.97	ng/ul	-0.01
19) Nitrobenzene-d5	8.90	128	28559	35.39	ng/ul	0.00
22) 2-Nitrophenol-d4	9.62	143	33356	36.81	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	10.15	165	57620	33.62	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.67	131	81192	41.45	ng/ul	-0.01
44) Dimethylphthalate-d6	13.80	166	202404	37.51	ng/ul	-0.01
47) Acenaphthylene-d8	14.08	160	231813	35.09	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.60	143	31419	32.55	ng/ul	-0.01
58) Fluorene-d10	15.39	176	170239	37.11	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.51	200	35139	32.87	ng/ul	-0.01
71) Anthracene-d10	17.25	188	281205	37.82	ng/ul	-0.01
79) Pyrene-d10	19.49	212	337151	38.96	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.39	264	383470	39.86	ng/ul	-0.01

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

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