

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030921\
 Data File : BP004945.D
 Acq On : 09 Mar 2021 15:54
 Operator : CG/JU
 Sample : PB134967BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK67

Quant Time: Mar 09 16:31:10 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP030921MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 09 15:13:30 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	32213	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	121859	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.39	164	80772	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.14	188	192234	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	234482	20.00	ng/ul	0.00
86) Perylene-d12	23.52	264	247983	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.23	96	4070	4.72	ng/uL	0.00
5) Phenol-d5	6.91	99	63311	25.02	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.07	67	35204	26.19	ng/ul	0.00
9) 2-Chlorophenol-d4	7.26	132	53218	26.62	ng/ul	0.00
13) 4-Methylphenol-d8	8.44	113	50824	23.84	ng/ul	0.00
19) Nitrobenzene-d5	8.89	128	25555	28.48	ng/ul	0.00
22) 2-Nitrophenol-d4	9.61	143	28332	28.13	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.15	165	49394	25.49	ng/ul	0.00
29) 4-Chloroaniline-d4	10.67	131	68419	32.26	ng/ul	0.00
44) Dimethylphthalate-d6	13.80	166	200319	33.70	ng/ul	0.00
47) Acenaphthylene-d8	14.07	160	210422	29.00	ng/ul	0.00
52) 4-Nitrophenol-d4	14.59	143	29739	29.12	ng/ul	0.00
58) Fluorene-d10	15.38	176	161606	31.63	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.50	200	36671	29.74	ng/ul	0.00
71) Anthracene-d10	17.24	188	286802	33.29	ng/ul	0.00
79) Pyrene-d10	19.48	212	363278	33.41	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	440700	34.75	ng/ul	0.00

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

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