

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP031921\
 Data File : BP005004.D
 Acq On : 19 Mar 2021 10:37
 Operator : CG/JU
 Sample : PB135014BL
 Misc : SFAM-MDL-S-BL02
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK014

Quant Time: Mar 19 11:08:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP031721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 19 09:59:18 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	43622	20.00	ng/ul	0.00
20) Naphthalene-d8	10.52	136	177381	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.37	164	121640	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.13	188	275552	20.00	ng/ul	0.00
79) Chrysene-d12	21.23	240	291322	20.00	ng/ul	0.00
88) Perylene-d12	23.52	264	279488	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	6675	6.05	ng/uL	0.00
4) Pyridine-d5	3.63	84	72567	26.95	ng/ul	0.00
7) Phenol-d5	6.90	99	110686	31.25	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.06	67	57748	30.92	ng/ul	0.00
11) 2-Chlorophenol-d4	7.26	132	95276	33.45	ng/ul	0.00
15) 4-Methylphenol-d8	8.43	113	89942	32.01	ng/ul	0.00
21) Nitrobenzene-d5	8.88	128	46345	34.18	ng/ul	0.00
24) 2-Nitrophenol-d4	9.60	143	54952	35.84	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.14	165	96011	32.56	ng/ul	0.00
31) 4-Chloroaniline-d4	10.66	131	129591	32.32	ng/ul	0.00
46) Dimethylphthalate-d6	13.79	166	328388	35.35	ng/ul	0.00
49) Acenaphthylene-d8	14.07	160	388151	35.78	ng/ul	0.00
54) 4-Nitrophenol-d4	14.59	143	47100	28.82	ng/ul	0.00
60) Fluorene-d10	15.37	176	277574	34.52	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.50	200	57569	29.93	ng/ul	0.00
73) Anthracene-d10	17.23	188	445088	34.34	ng/ul	0.00
81) Pyrene-d10	19.48	212	513592	35.48	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.37	264	553894	36.84	ng/ul	0.00

Target Compounds Ovalue

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

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