

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP031921\
 Data File : BP005002.D
 Acq On : 19 Mar 2021 09:29
 Operator : CG/JU
 Sample : PB135005BL
 Misc : SFAM-MDL-W-BL02
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK005

Quant Time: Mar 19 10:21:01 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	7.73	152	48812	20.00	ng/ul	0.01
20) Naphthalene-d8	10.52	136	187708	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.39	164	125482	20.00	ng/ul	0.01
64) Phenanthrene-d10	17.15	188	284658	20.00	ng/ul	0.00
79) Chrysene-d12	21.24	240	304068	20.00	ng/ul	0.00
88) Perylene-d12	23.53	264	297032	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	6844	5.54	ng/uL	0.00
4) Pyridine-d5	3.64	84	84992	28.21	ng/ul	0.00
7) Phenol-d5	6.90	99	122204	30.83	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.07	67	63490	30.38	ng/ul	0.00
11) 2-Chlorophenol-d4	7.26	132	106863	33.53	ng/ul	0.00
15) 4-Methylphenol-d8	8.44	113	99329	31.60	ng/ul	0.00
21) Nitrobenzene-d5	8.89	128	50187	34.97	ng/ul	0.01
24) 2-Nitrophenol-d4	9.61	143	58573	36.10	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.15	165	102458	32.83	ng/ul	0.00
31) 4-Chloroaniline-d4	10.67	131	137654	32.44	ng/ul	0.01
46) Dimethylphthalate-d6	13.80	166	337301	35.20	ng/ul	0.00
49) Acenaphthylene-d8	14.08	160	403285	36.04	ng/ul	0.01
54) 4-Nitrophenol-d4	14.59	143	49210	29.19	ng/ul	0.00
60) Fluorene-d10	15.39	176	287960	34.72	ng/ul	0.01
65) 4,6-Dinitro-2-methylphenol	15.51	200	57151	28.76	ng/ul	0.01
73) Anthracene-d10	17.25	188	458221	34.22	ng/ul	0.01
81) Pyrene-d10	19.49	212	536100	35.49	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.39	264	576165	36.06	ng/ul	0.01

Target Compounds Ovalue

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

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