

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060723\
 Data File : BP015397.D
 Acq On : 07 Jun 2023 14:12
 Operator : MA/JU
 Sample : PB153099BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK099

Quant Time: Jun 08 00:19:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 07 23:46:15 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.099	152	122239	20.000	ng/u1	0.00
20) Naphthalene-d8	10.928	136	537397	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.745	164	377302	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.498	188	901731	20.000	ng/u1	0.00
79) Chrysene-d12	21.580	240	871587	20.000	ng/u1	0.00
88) Perylene-d12	24.139	264	903230	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	25268	7.654	ng/uL	0.00
4) Pyridine-d5	3.852	84	302240	35.248	ng/u1	0.00
7) Phenol-d5	7.252	99	389619	34.598	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.422	67	262963	39.099	ng/u1	0.00
11) 2-Chlorophenol-d4	7.628	132	312773	38.307	ng/u1	0.00
15) 4-Methylphenol-d8	8.810	113	329709	35.673	ng/u1	0.00
21) Nitrobenzene-d5	9.275	128	168026	39.825	ng/u1	0.00
24) 2-Nitrophenol-d4	10.005	143	192509	39.955	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.546	165	312232	35.525	ng/u1	0.00
31) 4-Chloroaniline-d4	11.069	131	567085	44.950	ng/u1	0.00
46) Dimethylphthalate-d6	14.146	166	1310311	44.040	ng/u1	0.00
49) Acenaphthylene-d8	14.440	160	1341548	41.236	ng/u1	0.00
54) 4-Nitrophenol-d4	14.945	143	164027	30.786	ng/u1	0.00
60) Fluorene-d10	15.734	176	1038161	41.378	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200	188821	33.054	ng/u1	0.00
73) Anthracene-d10	17.598	188	1761293	42.930	ng/u1	0.00
81) Pyrene-d10	19.822	212	2125580	45.566	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.969	264	1985945	43.820	ng/u1	0.00

Target Compounds Qvalue

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

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