

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017555.D
 Acq On : 05 Oct 2023 07:53
 Operator : MA/JU
 Sample : PB156007BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK007

Quant Time: Oct 05 08:24:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 04 16:33:47 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	83256	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	334251	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	213015	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.422	188	493576	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	507894	20.000	ng/u1	0.00
88) Perylene-d12	24.133	264	575753	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	15524	7.455	ng/uL	0.00
4) Pyridine-d5	3.722	84	222277	40.778	ng/u1	0.00
7) Phenol-d5	7.093	99	281546	40.193	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	174473	40.818	ng/u1	0.00
11) 2-Chlorophenol-d4	7.457	132	226700	41.102	ng/u1	0.00
15) 4-Methylphenol-d8	8.657	113	225418	40.197	ng/u1	-0.01
21) Nitrobenzene-d5	9.146	128	106391	41.373	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	117339	40.457	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	216939	40.347	ng/u1	0.00
31) 4-Chloroaniline-d4	10.945	131	309006	40.342	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	734685	44.216	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	760082	41.485	ng/u1	0.00
54) 4-Nitrophenol-d4	14.875	143	110096	37.837	ng/u1	0.00
60) Fluorene-d10	15.633	176	620009	43.992	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	89750	29.167	ng/u1	0.00
73) Anthracene-d10	17.522	188	1026664	44.936	ng/u1	0.00
81) Pyrene-d10	19.774	212	1292062	45.782	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.962	264	1363106	46.913	ng/u1	-0.01

Target Compounds	Qvalue

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

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