

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092123\
 Data File : BP017257.D
 Acq On : 21 Sep 2023 09:11
 Operator : MA/JU
 Sample : PB155652BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK652

Quant Time: Sep 21 22:35:58 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 20 14:34:55 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	157293	20.000	ng/u1	0.00
20) Naphthalene-d8	10.763	136	656691	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.598	164	406727	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.369	188	911911	20.000	ng/u1	-0.01
79) Chrysene-d12	21.475	240	862811	20.000	ng/u1	-0.01
88) Perylene-d12	23.992	264	963742	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	24790	6.473	ng/uL	0.00
4) Pyridine-d5	3.740	84	348549	32.557	ng/u1	0.00
7) Phenol-d5	7.093	99	413203	31.985	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	253024	32.455	ng/u1	0.00
11) 2-Chlorophenol-d4	7.458	132	338662	33.053	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	332206	33.077	ng/u1	0.00
21) Nitrobenzene-d5	9.134	128	157365	33.147	ng/u1	0.00
24) 2-Nitrophenol-d4	9.852	143	179343	34.791	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	312110	32.047	ng/u1	0.00
31) 4-Chloroaniline-d4	10.928	131	463886	32.881	ng/u1	0.00
46) Dimethylphthalate-d6	14.016	166	1019385	34.986	ng/u1	0.00
49) Acenaphthylene-d8	14.298	160	1128838	33.686	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	147457	29.049	ng/u1	0.00
60) Fluorene-d10	15.598	176	874618	34.868	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.728	200	123051	23.941	ng/u1	0.00
73) Anthracene-d10	17.469	188	1436689	35.523	ng/u1	-0.01
81) Pyrene-d10	19.710	212	1725342	36.978	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.827	264	1713443	36.927	ng/u1	-0.02

Target Compounds Qvalue

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

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