

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017345.D
 Acq On : 26 Sep 2023 05:29
 Operator : MA/JU
 Sample : PB155717BL
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK717

Quant Time: Sep 26 06:02:22 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 26 04:52:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	203920	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	889584	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.587	164	563652	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.357	188	1287750	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	1169599	20.000	ng/u1	0.00
88) Perylene-d12	23.974	264	1270837	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	32903	6.627	ng/uL	0.00
4) Pyridine-d5	3.728	84	474403	34.181	ng/u1	0.00
7) Phenol-d5	7.081	99	571120	34.101	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	354169	35.041	ng/u1	0.00
11) 2-Chlorophenol-d4	7.446	132	462918	34.850	ng/u1	0.00
15) 4-Methylphenol-d8	8.640	113	458278	35.196	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	222159	34.544	ng/u1	0.00
24) 2-Nitrophenol-d4	9.840	143	247928	35.504	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	446999	33.881	ng/u1	0.00
31) 4-Chloroaniline-d4	10.916	131	656921	34.374	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	1483045	36.729	ng/u1	0.00
49) Acenaphthylene-d8	14.287	160	1592907	34.300	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	205208	29.171	ng/u1	0.00
60) Fluorene-d10	15.581	176	1260033	36.247	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	163771	22.564	ng/u1	0.00
73) Anthracene-d10	17.457	188	2046336	35.830	ng/u1	0.00
81) Pyrene-d10	19.698	212	2403362	37.999	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.810	264	2283100	37.313	ng/u1	0.00

Target Compounds Qvalue

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

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