

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092723\
 Data File : BP017397.D
 Acq On : 27 Sep 2023 14:53
 Operator : MA/JU
 Sample : PB155857BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK857

Quant Time: Sep 27 22:47:50 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	164966	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	757190	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.581	164	466977	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.351	188	931139	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	805486	20.000	ng/u1	0.00
88) Perylene-d12	23.974	264	891502	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	29254	7.284	ng/uL	0.00
4) Pyridine-d5	3.728	84	408539	36.386	ng/u1	0.00
7) Phenol-d5	7.075	99	489265	36.112	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	310891	38.022	ng/u1	0.00
11) 2-Chlorophenol-d4	7.440	132	395717	36.825	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	394243	37.428	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	193137	35.282	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	207637	34.933	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	375997	33.483	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	563267	34.627	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	1212977	36.259	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	1336364	34.733	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	129268	22.180	ng/u1	0.00
60) Fluorene-d10	15.581	176	1012661	35.162	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.710	200	108364	20.648	ng/u1	0.00
73) Anthracene-d10	17.451	188	1498685	36.291	ng/u1	0.00
81) Pyrene-d10	19.698	212	1713823	39.345	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.804	264	1610952	37.531	ng/u1	0.00

Target Compounds Qvalue

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

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