

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017150.D
 Acq On : 13 Sep 2023 15:41
 Operator : MA/JU
 Sample : PB155390BL
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK390

Quant Time: Sep 13 23:19:11 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 12 23:27:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	228635	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	942561	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.610	164	571601	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.381	188	1205530	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	831076	20.000	ng/ul	0.00
88) Perylene-d12	24.004	264	839475	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	32612	5.994	ng/uL	0.00
4) Pyridine-d5	3.746	84	372338	23.923	ng/ul	0.00
7) Phenol-d5	7.105	99	530154	28.517	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	341449	29.385	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	421736	28.436	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	418034	27.132	ng/ul	0.00
21) Nitrobenzene-d5	9.152	128	197539	29.930	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	215589	30.597	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	387236	28.417	ng/ul	0.00
31) 4-Chloroaniline-d4	10.940	131	352434	17.159	ng/ul	0.00
46) Dimethylphthalate-d6	14.028	166	1265093	31.102	ng/ul	0.00
49) Acenaphthylene-d8	14.310	160	1376845	29.831	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	166481	21.176	ng/ul	0.00
60) Fluorene-d10	15.610	176	1091647	31.038	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	148237	22.009	ng/ul	0.00
73) Anthracene-d10	17.481	188	1640653	32.031	ng/ul	0.00
81) Pyrene-d10	19.722	212	1707846	39.184	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.839	264	1267083	31.979	ng/ul	-0.01

Target Compounds Qvalue

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

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