

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017128.D
 Acq On : 12 Sep 2023 22:26
 Operator : MA/JU
 Sample : PB155340BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Sep 12 23:48:53 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	7.958	152	204176	20.000	ng/u1	0.00
20) Naphthalene-d8	10.781	136	846008	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.610	164	544709	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.381	188	1267300	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	1268722	20.000	ng/u1	0.00
88) Perylene-d12	24.004	264	1431306	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	29877	6.149	ng/uL	0.00
4) Pyridine-d5	3.752	84	399759	28.762	ng/u1	0.00
7) Phenol-d5	7.105	99	488505	29.424	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	305620	29.452	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	390561	29.489	ng/u1	0.00
15) 4-Methylphenol-d8	8.664	113	389811	28.331	ng/u1	0.00
21) Nitrobenzene-d5	9.152	128	181431	30.627	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	200536	31.709	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	365642	29.895	ng/u1	0.00
31) 4-Chloroaniline-d4	10.940	131	553444	30.021	ng/u1	0.00
46) Dimethylphthalate-d6	14.028	166	1251094	32.276	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	1335704	30.368	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	199023	26.566	ng/u1	0.00
60) Fluorene-d10	15.610	176	1069344	31.905	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200	160436	22.659	ng/u1	-0.01
73) Anthracene-d10	17.481	188	1778002	33.021	ng/u1	0.00
81) Pyrene-d10	19.722	212	2240316	33.670	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.839	264	2269133	33.589	ng/u1	-0.01

Target Compounds Qvalue

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

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