

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017185.D
 Acq On : 14 Sep 2023 19:49
 Operator : MA/JU
 Sample : PB155424BL
 Misc : GCMS CONFIRMATION
 ALS Vial : 74 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK424

Quant Time: Sep 14 22:49:59 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	7.946	152	231561	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.775	136	887021	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.610	164	523684	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.380	188	1256119	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	973862	20.000	ng/u1	0.00
88) Perylene-d12	24.004	264	874705	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	119	0.022	ng/uL	-0.01
4) Pyridine-d5	3.758	84	35	0.002	ng/u1	0.00
7) Phenol-d5	7.093	99	14	0.001	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.293	67	9	0.001	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	17	0.001	ng/u1	-0.01
15) 4-Methylphenol-d8	8.657	113	20	0.001	ng/u1	-0.01
21) Nitrobenzene-d5	9.140	128	45	0.007	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.869	143	23	0.003	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	15	0.001	ng/u1	-0.10
31) 4-Chloroaniline-d4	10.934	131	61	0.003	ng/u1	-0.01
46) Dimethylphthalate-d6	14.039	166	96	0.003	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	307	0.007	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	27	0.004	ng/u1	0.00
60) Fluorene-d10	15.610	176	633	0.020	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.728	200	13	0.002	ng/u1	-0.02
73) Anthracene-d10	17.474	188	4039	0.076	ng/u1	-0.01
81) Pyrene-d10	19.716	212	1322	0.026	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.004	264	654118	15.844	ng/u1	0.15

Target Compounds Qvalue

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

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