

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
 Data File : BP017997.D
 Acq On : 09 Nov 2023 20:07
 Operator : MA/JU
 Sample : PB156742BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK742

Quant Time: Nov 09 23:26:11 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 09 22:34:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	62054	20.000	ng/u1	0.00
20) Naphthalene-d8	10.687	136	258960	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.539	164	168937	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.322	188	398052	20.000	ng/u1	0.00
79) Chrysene-d12	21.445	240	384751	20.000	ng/u1	0.00
88) Perylene-d12	23.957	264	420807	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	10566	5.829	ng/uL	0.00
4) Pyridine-d5	3.681	84	132799	28.187	ng/u1	0.00
7) Phenol-d5	7.028	99	157512	27.010	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	102698	29.072	ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	130160	27.896	ng/u1	0.00
15) 4-Methylphenol-d8	8.587	113	129628	27.714	ng/u1	0.00
21) Nitrobenzene-d5	9.069	128	63312	27.301	ng/u1	0.00
24) 2-Nitrophenol-d4	9.787	143	72673	28.158	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	112938	24.427	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	167702	26.468	ng/u1	0.00
46) Dimethylphthalate-d6	13.963	166	472750	30.441	ng/u1	0.00
49) Acenaphthylene-d8	14.239	160	468795	28.157	ng/u1	0.00
54) 4-Nitrophenol-d4	14.804	143	51191	22.506	ng/u1	0.00
60) Fluorene-d10	15.545	176	374506	29.095	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	66741	24.013	ng/u1	0.00
73) Anthracene-d10	17.422	188	617370	28.946	ng/u1	0.00
81) Pyrene-d10	19.674	212	753037	30.411	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.792	264	720883	29.473	ng/u1	0.00

Target Compounds Qvalue

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

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