

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
 Data File : BP017502.D
 Acq On : 03 Oct 2023 05:53
 Operator : MA/JU
 Sample : 04588-04
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHS6

Quant Time: Oct 03 06:30:00 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 30 04:02:20 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	123950	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.775	136	509889	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.634	164	299789	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.422	188	658571	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	648587	20.000	ng/u1	0.00
88) Perylene-d12	24.139	264	724484	20.000	ng/u1	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	15446	5.118	ng/uL	0.00
4) Pyridine-d5	3.722	84	138736	16.445	ng/u1	0.00
7) Phenol-d5	7.099	99	99493	9.773	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	282818	46.035	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.458	132	291764	36.136	ng/u1	-0.01
15) 4-Methylphenol-d8	8.657	113	181690	22.957	ng/u1	-0.01
21) Nitrobenzene-d5	9.146	128	185792	50.402	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.857	143	199374	49.812	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.393	165	329050	43.514	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.946	131	410377	37.463	ng/u1	-0.01
46) Dimethylphthalate-d6	14.045	166	1085869	50.562	ng/u1	-0.01
49) Acenaphthylene-d8	14.328	160	1223242	49.524	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.898	143	24255	6.483	ng/u1	0.02
60) Fluorene-d10	15.634	176	942203	50.961	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.775	200	120681	32.512	ng/u1	-0.01
73) Anthracene-d10	17.522	188	1573160	53.860	ng/u1	0.00
81) Pyrene-d10	19.780	212	1956137	55.772	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.974	264	1909331	54.737	ng/u1	0.02

Target Compounds	Qvalue

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

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