

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006932.D
 Acq On : 10 Sep 2021 10:03
 Operator : CG/JU
 Sample : M3552-04 10X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AH3

Quant Time: Sep 10 10:53:50 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 09 11:07:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	298060	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	1216618	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.557	164	754223	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.304	188	1531563	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.392	240	1426257	20.000	ng/u1	0.00
88) Perylene-d12	23.821	264	1461250	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	22066	2.946	ng/uL	0.01
4) Pyridine-d5	3.799	84	174512	8.941	ng/u1	0.02
7) Phenol-d5	7.093	99	165994	7.577	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	491453	39.062	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	495214	28.002	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	306299	17.805	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	288849	39.197	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	273408	37.685	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.351	165	523110	30.579	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	650958	27.516	ng/u1	0.00
46) Dimethylphthalate-d6	13.969	166	1854189	37.957	ng/u1	0.00
49) Acenaphthylene-d8	14.251	160	2269475	36.682	ng/u1	0.00
54) 4-Nitrophenol-d4	14.739	143	39970	4.710	ng/u1	0.00
60) Fluorene-d10	15.551	176	1680207	38.972	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	209907	30.981	ng/u1	0.00
73) Anthracene-d10	17.404	188	2850717	44.959	ng/u1	-0.01
81) Pyrene-d10	19.639	212	3247951	48.625	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.668	264	3020172	42.843	ng/u1	-0.01

Target Compounds	Qvalue

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

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