

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006901.D
 Acq On : 09 Sep 2021 15:32
 Operator : CG/JU
 Sample : M3552-04 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AH3

Quant Time: Sep 09 16:43:03 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	477454	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	2131612	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.563	164	1349536	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.310	188	2850761	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.398	240	2902070	20.000	ng/u1	# 0.00
88) Perylene-d12	23.839	264	3059804	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	30997	2.584	ng/uL	0.00
4) Pyridine-d5	3.787	84	273744	8.755	ng/u1	0.00
7) Phenol-d5	7.093	99	336661	9.594	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	974594	48.358	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	961916	33.955	ng/u1	0.00
15) 4-Methylphenol-d8	8.640	113	674884	24.490	ng/u1	0.00
21) Nitrobenzene-d5	9.099	128	676778	52.417	ng/u1	0.00
24) 2-Nitrophenol-d4	9.822	143	681740	53.632	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.358	165	1178880	39.332	ng/u1	0.00
31) 4-Chloroaniline-d4	10.875	131	1419983	34.258	ng/u1	0.00
46) Dimethylphthalate-d6	13.975	166	4149798	47.477	ng/u1	0.00
49) Acenaphthylene-d8	14.257	160	5094321	46.019	ng/u1	0.00
54) 4-Nitrophenol-d4	14.746	143	134666	8.868	ng/u1	0.00
60) Fluorene-d10	15.557	176	3679643	47.699	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	667468	52.927	ng/u1	0.00
73) Anthracene-d10	17.416	188	6201696	52.547	ng/u1	0.00
81) Pyrene-d10	19.645	212	7247796	53.327	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.686	264	7578734	51.342	ng/u1	0.00

Target Compounds Qvalue

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

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