

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111022\
 Data File : BP012602.D
 Acq On : 10 Nov 2022 16:41
 Operator : CG/JU
 Sample : N5493-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0AA6

Quant Time: Nov 10 21:35:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 09 23:28:28 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.757	152	351397	20.000	ng/u1	0.00
20) Naphthalene-d8	10.557	136	1590377	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.416	164	1090932	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.180	188	2453820	20.000	ng/u1	0.00
79) Chrysene-d12	21.280	240	2723716	20.000	ng/u1	0.00
88) Perylene-d12	23.657	264	2752416	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	29810	3.458	ng/uL	0.00
4) Pyridine-d5	3.652	84	167918	6.892	ng/u1	0.01
7) Phenol-d5	6.952	99	172403	5.820	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	473794	26.732	ng/u1	0.00
11) 2-Chlorophenol-d4	7.299	132	496676	21.788	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	327995	13.997	ng/u1	0.00
21) Nitrobenzene-d5	8.934	128	330613	28.095	ng/u1	0.00
24) 2-Nitrophenol-d4	9.651	143	353173	27.326	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	590943	25.379	ng/u1	0.00
31) 4-Chloroaniline-d4	10.716	131	750864	22.893	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	2372324	31.549	ng/u1	0.00
49) Acenaphthylene-d8	14.104	160	2490995	29.283	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	60963	5.149	ng/u1	0.02
60) Fluorene-d10	15.416	176	2098223	32.909	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.551	200	379285	27.946	ng/u1	0.00
73) Anthracene-d10	17.281	188	3587507	33.666	ng/u1	0.00
81) Pyrene-d10	19.522	212	4328594	28.359	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.504	264	4524271	33.344	ng/u1	0.00

Target Compounds Qvalue

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

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