

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP012020.D
 Acq On : 07 Oct 2022 21:58
 Operator : CG/JU
 Sample : N4951-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8T5

Quant Time: Oct 07 23:11:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 07 03:22:44 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	202034	20.000	ng/ul	0.00
20) Naphthalene-d8	10.681	136	858120	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	492723	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1058407	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1027040	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1046291	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	15794	3.448	ng/uL	0.00
4) Pyridine-d5	3.728	84	136595	10.180	ng/ul	0.01
7) Phenol-d5	7.040	99	128279	7.355	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	324218	33.536	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132	362108	26.300	ng/ul	0.00
15) 4-Methylphenol-d8	8.587	113	220102	15.266	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	221590	33.902	ng/ul	0.00
24) 2-Nitrophenol-d4	9.763	143	220008	31.461	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	351204	27.152	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	450891	22.866	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	1176804	33.399	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1462474	36.409	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	29948	4.104	ng/ul	0.02
60) Fluorene-d10	15.516	176	1100991	35.981	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	119679	18.682	ng/ul	0.00
73) Anthracene-d10	17.381	188	1804406	37.426	ng/ul	0.00
81) Pyrene-d10	19.616	212	2028037	38.990	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	2012369	38.421	ng/ul	0.00

Target Compounds	Qvalue

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

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