

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011953.D
 Acq On : 06 Oct 2022 04:48
 Operator : CG/JU
 Sample : N4956-01
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8Q7

Quant Time: Oct 06 05:29:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 05 23:18:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	297398	20.000	ng/u1	0.00
20) Naphthalene-d8	10.687	136	1269075	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	771515	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.281	188	1632868	20.000	ng/u1	0.00
79) Chrysene-d12	21.369	240	1552772	20.000	ng/u1	0.00
88) Perylene-d12	23.798	264	1580314	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	29622	4.393	ng/uL	0.00
4) Pyridine-d5	3.734	84	60654	3.071	ng/u1	0.01
7) Phenol-d5	7.046	99	126979	4.946	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	367802	25.845	ng/u1	0.00
11) 2-Chlorophenol-d4	7.411	132	402146	19.842	ng/u1	0.00
15) 4-Methylphenol-d8	8.587	113	242395	11.421	ng/u1	-0.01
21) Nitrobenzene-d5	9.046	128	253841	26.260	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	248929	24.070	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	433446	22.659	ng/u1	0.00
31) 4-Chloroaniline-d4	10.828	131	549841	18.854	ng/u1	0.00
46) Dimethylphthalate-d6	13.934	166	1517941	27.513	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	1732435	27.545	ng/u1	0.00
54) 4-Nitrophenol-d4	14.740	143	34144	2.989	ng/u1	0.01
60) Fluorene-d10	15.522	176	1388725	28.984	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	189009	19.124	ng/u1	0.00
73) Anthracene-d10	17.381	188	2291606	30.809	ng/u1	0.00
81) Pyrene-d10	19.616	212	2622697	33.351	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.639	264	2504291	31.656	ng/u1	0.00

Target Compounds Qvalue

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

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