

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011895.D
 Acq On : 04 Oct 2022 04:44
 Operator : CG/JU
 Sample : N4859-11
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E10051

Quant Time: Oct 04 05:45:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 01 00:53:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	263930	20.000	ng/u1	0.00
20) Naphthalene-d8	10.693	136	1192327	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.534	164	700991	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.286	188	1483688	20.000	ng/u1	-0.01
79) Chrysene-d12	21.374	240	1715927	20.000	ng/u1	-0.01
88) Perylene-d12	23.810	264	1840161	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	34948	5.840	ng/uL	0.00
4) Pyridine-d5	3.728	84	126877	7.238	ng/u1	0.00
7) Phenol-d5	7.052	99	127218	5.583	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	374486	29.651	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.416	132	417541	23.214	ng/u1	0.00
15) 4-Methylphenol-d8	8.599	113	265961	14.120	ng/u1	0.00
21) Nitrobenzene-d5	9.052	128	291466	32.093	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.775	143	294598	30.319	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.316	165	497628	27.689	ng/u1	0.00
31) 4-Chloroaniline-d4	10.834	131	654954	23.904	ng/u1	-0.01
46) Dimethylphthalate-d6	13.945	166	1607504	32.068	ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	1877001	32.846	ng/u1	0.00
54) 4-Nitrophenol-d4	14.751	143	36910	3.556	ng/u1	0.01
60) Fluorene-d10	15.528	176	1496816	34.383	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	231431	25.771	ng/u1	0.00
73) Anthracene-d10	17.386	188	2569203	38.014	ng/u1	-0.01
81) Pyrene-d10	19.622	212	3254209	37.447	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.651	264	3684829	40.002	ng/u1	-0.01

Target Compounds	Qvalue

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

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