

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011879.D
 Acq On : 03 Oct 2022 19:06
 Operator : CG/JU
 Sample : N4859-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E10018

Quant Time: Oct 03 23:33:03 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	7.893	152	233015	20.000	ng/u1	0.00
20) Naphthalene-d8	10.698	136	929513	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.534	164	532211	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.292	188	1100830	20.000	ng/u1	0.00
79) Chrysene-d12	21.374	240	1171479	20.000	ng/u1	-0.01
88) Perylene-d12	23.810	264	1230047	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	28478	5.390	ng/uL	0.00
4) Pyridine-d5	3.734	84	110355	7.131	ng/u1	0.00
7) Phenol-d5	7.052	99	110963	5.516	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	301624	27.051	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.416	132	345223	21.740	ng/u1	0.00
15) 4-Methylphenol-d8	8.599	113	204800	12.316	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	212306	29.987	ng/u1	0.00
24) 2-Nitrophenol-d4	9.781	143	211567	27.930	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	364168	25.992	ng/u1	0.00
31) 4-Chloroaniline-d4	10.840	131	443153	20.747	ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	1177651	30.943	ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	1334887	30.767	ng/u1	0.00
54) 4-Nitrophenol-d4	14.739	143	37788	4.795	ng/u1	0.00
60) Fluorene-d10	15.528	176	1064000	32.192	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	159657	23.962	ng/u1	0.00
73) Anthracene-d10	17.392	188	1792592	35.748	ng/u1	0.00
81) Pyrene-d10	19.622	212	2256124	38.027	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.651	264	2370092	38.491	ng/u1	-0.01

Target Compounds Qvalue

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

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