

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011873.D
 Acq On : 03 Oct 2022 15:41
 Operator : CG/JU
 Sample : N4749-08
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8Q3

Quant Time: Oct 03 23:31:50 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	216967	20.000	ng/u1	0.00
20) Naphthalene-d8	10.699	136	856960	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.534	164	488369	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.292	188	1021603	20.000	ng/u1	0.00
79) Chrysene-d12	21.374	240	1183250	20.000	ng/u1	-0.01
88) Perylene-d12	23.810	264	1232378	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	23922	4.863	ng/uL	0.00
4) Pyridine-d5	3.734	84	121382	8.423	ng/u1	0.00
7) Phenol-d5	7.052	99	128518	6.861	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	336339	32.395	ng/u1	0.00
11) 2-Chlorophenol-d4	7.422	132	375938	25.425	ng/u1	0.00
15) 4-Methylphenol-d8	8.605	113	247307	15.972	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	235473	36.075	ng/u1	0.00
24) 2-Nitrophenol-d4	9.781	143	225262	32.256	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	402214	31.138	ng/u1	0.00
31) 4-Chloroaniline-d4	10.840	131	352352	17.893	ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	1411889	40.428	ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	1586413	39.847	ng/u1	0.00
54) 4-Nitrophenol-d4	14.745	143	40203	5.559	ng/u1	0.00
60) Fluorene-d10	15.528	176	1258795	41.505	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	165772	26.809	ng/u1	0.00
73) Anthracene-d10	17.392	188	2130546	45.783	ng/u1	0.00
81) Pyrene-d10	19.627	212	2679910	44.721	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.657	264	2942247	47.693	ng/u1	0.00

Target Compounds	Qvalue

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

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