

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011893.D
 Acq On : 04 Oct 2022 03:36
 Operator : CG/JU
 Sample : N4749-05
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8N4

Quant Time: Oct 04 04:37:19 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	265311	20.000 ng/u1	0.00
20) Naphthalene-d8	10.699	136	1087317	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.534	164	651303	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.287	188	1378752	20.000 ng/u1	-0.01
79) Chrysene-d12	21.375	240	1438507	20.000 ng/u1	-0.01
88) Perylene-d12	23.810	264	1490255	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.299	96	31829	5.291 ng/uL	0.00
4) Pyridine-d5	3.728	84	135604	7.696 ng/u1	0.00
7) Phenol-d5	7.052	99	139388	6.086 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	352188	27.741 ng/u1	-0.01
11) 2-Chlorophenol-d4	7.416	132	401775	22.221 ng/u1	0.00
15) 4-Methylphenol-d8	8.599	113	259456	13.703 ng/u1	0.00
21) Nitrobenzene-d5	9.052	128	241412	29.149 ng/u1	-0.01
24) 2-Nitrophenol-d4	9.775	143	241768	27.285 ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.316	165	434945	26.538 ng/u1	0.00
31) 4-Chloroaniline-d4	10.834	131	489996	19.611 ng/u1	-0.01
46) Dimethylphthalate-d6	13.946	166	1553671	33.358 ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	1700396	32.025 ng/u1	0.00
54) 4-Nitrophenol-d4	14.740	143	46692	4.841 ng/u1	0.00
60) Fluorene-d10	15.528	176	1390301	34.373 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	205288	24.600 ng/u1	0.00
73) Anthracene-d10	17.387	188	2299653	36.616 ng/u1	-0.01
81) Pyrene-d10	19.622	212	2883199	39.576 ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.651	264	2908880	38.993 ng/u1	-0.01

Target Compounds Qvalue

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

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