

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011975.D
 Acq On : 06 Oct 2022 18:00
 Operator : CG/JU
 Sample : N4923-10
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E10016

Quant Time: Oct 06 23:07:59 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	358668	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	1398727	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	718238	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.280	188	1330870	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1467020	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1573893	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	32011	3.936	ng/uL	0.00
4) Pyridine-d5	3.734	84	126759	5.321	ng/ul	0.01
7) Phenol-d5	7.046	99	143031	4.619	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	415746	24.223	ng/ul	0.00
11) 2-Chlorophenol-d4	7.410	132	431213	17.642	ng/ul	0.00
15) 4-Methylphenol-d8	8.593	113	272384	10.642	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	281436	26.416	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	283118	24.838	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	432693	20.523	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	462692	14.395	ng/ul	0.00
46) Dimethylphthalate-d6	13.939	166	1334395	25.980	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	1580715	26.997	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	26319	2.475	ng/ul	0.02
60) Fluorene-d10	15.522	176	1190376	26.687	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	152889	18.980	ng/ul	0.00
73) Anthracene-d10	17.380	188	1729133	28.522	ng/ul	0.00
81) Pyrene-d10	19.616	212	2081887	28.021	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	2267147	28.776	ng/ul	0.00

Target Compounds Qvalue

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

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