

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011962.D
 Acq On : 06 Oct 2022 10:29
 Operator : CG/JU
 Sample : N4956-06
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 06 23:04:29 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	7.875	152	248586	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	1046679	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	636925	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1351649	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1303868	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1293598	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	24492	4.345	ng/uL	0.00
4) Pyridine-d5	3.734	84	51622	3.127	ng/ul	0.01
7) Phenol-d5	7.040	99	115881	5.400	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	320763	26.965	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132	367793	21.710	ng/ul	0.00
15) 4-Methylphenol-d8	8.587	113	221283	12.473	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	225951	28.342	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	221880	26.013	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	414568	26.277	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	432290	17.973	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	1392586	30.575	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1544927	29.754	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	31572	3.347	ng/ul	0.00
60) Fluorene-d10	15.516	176	1274588	32.223	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	136476	16.682	ng/ul	0.00
73) Anthracene-d10	17.380	188	2099346	34.097	ng/ul	0.00
81) Pyrene-d10	19.616	212	2458865	37.237	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	2321700	35.853	ng/ul	0.00

Target Compounds Qvalue

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

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