

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012121.D
 Acq On : 13 Oct 2022 15:19
 Operator : CG/JU
 Sample : N4982-09
 Misc :
 ALS Vial : 92 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 13 23:33:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	280014	20.000	ng/ul	0.00
20) Naphthalene-d8	10.657	136	1123977	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.498	164	615361	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.257	188	1262649	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	1367357	20.000	ng/ul	0.00
88) Perylene-d12	23.768	264	1483419	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	31761	4.472	ng/uL	0.00
4) Pyridine-d5	3.705	84	132833	7.033	ng/ul	0.00
7) Phenol-d5	7.016	99	168842	7.497	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	428840	33.548	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	487057	27.246	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	283809	16.407	ng/ul	0.00
21) Nitrobenzene-d5	9.016	128	298582	37.614	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	296592	36.848	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	481723	30.881	ng/ul	0.00
31) 4-Chloroaniline-d4	10.798	131	610021	26.537	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	1399255	35.904	ng/ul	0.00
49) Acenaphthylene-d8	14.192	160	1705636	35.393	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	40968	5.355	ng/ul	0.00
60) Fluorene-d10	15.492	176	1290861	36.610	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	184434	27.663	ng/ul	0.00
73) Anthracene-d10	17.357	188	2041515	37.234	ng/ul	0.00
81) Pyrene-d10	19.598	212	2528704	36.098	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.609	264	2719116	37.707	ng/ul	0.00

Target Compounds Qvalue

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

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