

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012440.D
 Acq On : 01 Nov 2022 22:28
 Operator : CG/JU
 Sample : N5300-06
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA88

Quant Time: Nov 02 00:53:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 01 01:03:58 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	295164	20.000	ng/u1	0.00
20) Naphthalene-d8	10.587	136	1338619	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.439	164	893082	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.204	188	1979066	20.000	ng/u1	0.00
79) Chrysene-d12	21.298	240	1880003	20.000	ng/u1	0.00
88) Perylene-d12	23.686	264	1596554	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	30669	4.184	ng/uL	0.00
4) Pyridine-d5	3.663	84	171013	8.121	ng/u1	0.00
7) Phenol-d5	6.963	99	181189	6.844	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	482228	30.513	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	488382	25.262	ng/u1	0.00
15) 4-Methylphenol-d8	8.504	113	339872	16.232	ng/u1	0.00
21) Nitrobenzene-d5	8.957	128	344178	39.649	ng/u1	0.00
24) 2-Nitrophenol-d4	9.681	143	373956	41.505	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	602913	30.877	ng/u1	0.00
31) 4-Chloroaniline-d4	10.739	131	812410	27.943	ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	2162355	35.369	ng/u1	0.00
49) Acenaphthylene-d8	14.133	160	2440088	34.925	ng/u1	0.00
54) 4-Nitrophenol-d4	14.675	143	64728	5.594	ng/u1	0.00
60) Fluorene-d10	15.439	176	1905675	36.407	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	200609	19.469	ng/u1	0.00
73) Anthracene-d10	17.304	188	3230989	37.890	ng/u1	0.00
81) Pyrene-d10	19.545	212	3745919	39.686	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.527	264	3030231	38.844	ng/u1	-0.02

Target Compounds Qvalue

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

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