

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013243.D
 Acq On : 22 Dec 2022 13:22
 Operator : CG/JU
 Sample : N6015-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 22 21:46:38 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 21 23:38:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	653162	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	2750572	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	1709617	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	3391137	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	2234378	20.000	ng/ul	0.00
88) Perylene-d12	23.910	264	2557185	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	70443	4.611	ng/uL	0.00
4) Pyridine-d5	3.764	84	829817	19.122	ng/ul	0.00
7) Phenol-d5	7.105	99	1356735	25.502	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	847238	26.399	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	1108738	26.474	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	976847	22.409	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	539450	26.693	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	615933	27.072	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	1155366	26.144	ng/ul	0.00
31) 4-Chloroaniline-d4	10.899	131	844833	13.542	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	3336591	25.394	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	3813012	26.410	ng/ul	0.00
54) 4-Nitrophenol-d4	14.787	143	395755	16.745	ng/ul	0.00
60) Fluorene-d10	15.581	176	2903391	26.119	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	407522	18.674	ng/ul	0.00
73) Anthracene-d10	17.445	188	3986810	26.059	ng/ul	0.00
81) Pyrene-d10	19.675	212	3895746	30.375	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.745	264	3375081	26.479	ng/ul	0.00

Target Compounds Qvalue

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

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