

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013313.D
 Acq On : 24 Dec 2022 23:09
 Operator : CG/JU
 Sample : N6016-17
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 25 00:25:36 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 24 00:19:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	632350	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	2382858	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.586	164	1234995	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	2512962	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	2746807	20.000	ng/ul	0.00
88) Perylene-d12	23.915	264	3212682	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	49273	3.332	ng/uL	0.00
4) Pyridine-d5	3.775	84	57507	1.369	ng/ul	0.02
7) Phenol-d5	7.099	99	868371	16.859	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.263	67	601116	19.347	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	748929	18.471	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	492467	11.669	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	353133	20.170	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	384982	19.532	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	661144	17.269	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	545250	10.088	ng/ul	0.00
46) Dimethylphthalate-d6	13.987	166	1793500	18.896	ng/ul	-0.01
49) Acenaphthylene-d8	14.275	160	1995128	19.130	ng/ul	0.00
54) 4-Nitrophenol-d4	14.787	143	184147	10.786	ng/ul	0.00
60) Fluorene-d10	15.581	176	1481938	18.455	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	121799	7.532	ng/ul	0.00
73) Anthracene-d10	17.439	188	2151860	18.981	ng/ul	0.00
81) Pyrene-d10	19.674	212	2998840	19.020	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	3171333	19.804	ng/ul	0.00

Target Compounds Qvalue

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

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