

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013152.D
 Acq On : 20 Dec 2022 06:50
 Operator : CG/JU
 Sample : N6003-01
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXX0

Quant Time: Dec 20 21:43:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 19 21:43:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	498299	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	1977810	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.598	164	1132123	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.351	188	2127931	20.000	ng/ul	0.00
79) Chrysene-d12	21.439	240	1851960	20.000	ng/ul	0.00
88) Perylene-d12	23.916	264	2329384	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	64224	5.511	ng/uL	0.00
4) Pyridine-d5	3.770	84	301332	9.102	ng/ul	0.00
7) Phenol-d5	7.111	99	1221442	30.094	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	811896	33.160	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	1014960	31.766	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	947284	28.484	ng/ul	0.00
21) Nitrobenzene-d5	9.122	128	503542	34.652	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	569102	34.787	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	983120	30.939	ng/ul	0.00
31) 4-Chloroaniline-d4	10.905	131	1062158	23.677	ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	2962276	34.046	ng/ul	0.00
49) Acenaphthylene-d8	14.293	160	3260621	34.104	ng/ul	0.00
54) 4-Nitrophenol-d4	14.793	143	374519	23.929	ng/ul	0.00
60) Fluorene-d10	15.587	176	2487661	33.795	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	380685	27.800	ng/ul	0.00
73) Anthracene-d10	17.451	188	3429927	35.728	ng/ul	0.00
81) Pyrene-d10	19.681	212	3715762	34.955	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.757	264	3979785	34.277	ng/ul	0.00

Target Compounds Qvalue

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

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