

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012471.D
 Acq On : 03 Nov 2022 02:54
 Operator : CG/JU
 Sample : N5321-18
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBWR7

Quant Time: Nov 03 03:48:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 03 00:09:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	294671	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	1373824	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	939191	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	2111501	20.000	ng/u1	0.00
79) Chrysene-d12	21.292	240	2035032	20.000	ng/u1	0.00
88) Perylene-d12	23.674	264	1828584	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	32252	4.407	ng/uL	0.00
4) Pyridine-d5	3.658	84	204146	9.711	ng/u1	0.00
7) Phenol-d5	6.963	99	226085	8.554	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	522239	33.100	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	557825	28.902	ng/u1	0.00
15) 4-Methylphenol-d8	8.504	113	428798	20.513	ng/u1	0.00
21) Nitrobenzene-d5	8.951	128	361863	40.618	ng/u1	0.00
24) 2-Nitrophenol-d4	9.675	143	396064	42.832	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	708681	35.364	ng/u1	0.00
31) 4-Chloroaniline-d4	10.734	131	922082	30.903	ng/u1	0.00
46) Dimethylphthalate-d6	13.851	166	2537267	39.464	ng/u1	0.00
49) Acenaphthylene-d8	14.128	160	2722176	37.050	ng/u1	0.00
54) 4-Nitrophenol-d4	14.675	143	91334	7.506	ng/u1	0.01
60) Fluorene-d10	15.433	176	2155112	39.151	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	399998	36.385	ng/u1	0.00
73) Anthracene-d10	17.298	188	3584118	39.395	ng/u1	0.00
81) Pyrene-d10	19.539	212	4190574	41.015	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.521	264	3630158	40.630	ng/u1	0.00

Target Compounds Qvalue

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

