

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012105.D
 Acq On : 13 Oct 2022 05:33
 Operator : CG/JU
 Sample : N5026-06
 Misc :
 ALS Vial : 75 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BX4

Quant Time: Oct 13 06:15:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	236417	20.000	ng/u1	0.00
20) Naphthalene-d8	10.657	136	968234	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.498	164	578817	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	1283483	20.000	ng/u1	0.00
79) Chrysene-d12	21.345	240	1331664	20.000	ng/u1	0.00
88) Perylene-d12	23.763	264	1427408	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	24271	4.048	ng/uL	0.00
4) Pyridine-d5	3.705	84	114307	7.168	ng/u1	0.00
7) Phenol-d5	7.022	99	144889	7.620	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	338028	31.320	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	400896	26.562	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	251683	17.233	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	241415	35.304	ng/u1	0.00
24) 2-Nitrophenol-d4	9.740	143	245690	35.434	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	418429	31.138	ng/u1	0.00
31) 4-Chloroaniline-d4	10.799	131	443910	22.417	ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	1412290	38.527	ng/u1	0.00
49) Acenaphthylene-d8	14.193	160	1624210	35.831	ng/u1	0.00
54) 4-Nitrophenol-d4	14.716	143	46933	6.522	ng/u1	0.00
60) Fluorene-d10	15.493	176	1318071	39.742	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	206409	30.456	ng/u1	0.00
73) Anthracene-d10	17.357	188	2265100	40.641	ng/u1	0.00
81) Pyrene-d10	19.592	212	2748158	40.283	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.610	264	2855049	41.145	ng/u1	0.00

Target Compounds Qvalue

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

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